

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
 Data File : BP022106.D
 Acq On : 28 Sep 2024 15:46
 Operator : RC/JU
 Sample : P3910-10DL 10X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DDC29DL

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022106.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.757	464	471	479	rVV2	1212980	1815391	2.24%	0.813%
2	6.287	556	561	566	rBV	338683	480035	0.59%	0.215%
3	6.710	623	633	638	rBV3	496610	1139931	1.40%	0.511%
4	6.769	638	643	646	rVV	567270	800150	0.99%	0.359%
5	6.804	646	649	654	rVV	705711	1022434	1.26%	0.458%
6	6.869	654	660	666	rVV3	858915	1619041	2.00%	0.725%
7	6.940	666	672	676	rVV	611717	895538	1.10%	0.401%
8	6.987	676	680	685	rVB	361571	519426	0.64%	0.233%
9	7.098	693	699	702	rVV	520493	824139	1.02%	0.369%
10	7.134	702	705	709	rVV	411428	624488	0.77%	0.280%
11	7.228	713	721	725	rVV2	467102	883042	1.09%	0.396%
12	7.334	733	739	748	rVV2	2793225	5489663	6.77%	2.460%
13	7.622	777	788	793	rVV4	179219	575266	0.71%	0.258%
14	7.681	793	798	805	rVV2	657350	1366197	1.68%	0.612%
15	7.769	805	813	817	rVV	1371609	2365958	2.92%	1.060%
16	7.816	817	821	828	rVB2	429775	773555	0.95%	0.347%
17	7.916	835	838	842	rVV	508211	741001	0.91%	0.332%
18	7.987	847	850	857	rVB3	282427	481988	0.59%	0.216%
19	8.128	868	874	878	rBV	322109	581202	0.72%	0.260%
20	8.228	887	891	897	rVB2	436544	845738	1.04%	0.379%
21	8.328	904	908	911	rBV2	645036	1144662	1.41%	0.513%
22	8.457	924	930	933	rBV	540769	838997	1.03%	0.376%
23	8.492	933	936	944	rVB	401321	729664	0.90%	0.327%
24	8.640	956	961	965	rBV	363228	561629	0.69%	0.252%
25	8.687	965	969	977	rVB	426351	709986	0.88%	0.318%
26	8.792	977	987	995	rBV2	536383	1301009	1.60%	0.583%
27	8.910	997	1007	1013	rVB	2535624	4092970	5.04%	1.834%
28	9.034	1018	1028	1035	rVB8	325101	920244	1.13%	0.412%
29	9.116	1035	1042	1046	rBV2	615852	1131327	1.39%	0.507%
30	9.557	1108	1117	1120	rBV2	224012	461958	0.57%	0.207%
31	9.604	1120	1125	1129	rVV3	400829	832860	1.03%	0.373%
32	9.645	1130	1132	1142	rVV4	231826	713984	0.88%	0.320%
33	9.751	1147	1150	1155	rVV	369489	573819	0.71%	0.257%
34	9.822	1155	1162	1165	rVV3	278324	492884	0.61%	0.221%
35	9.892	1165	1174	1178	rVV4	338023	1001896	1.23%	0.449%
36	10.163	1214	1220	1224	rBV2	311562	519123	0.64%	0.233%

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Integrator: RTE

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Title : SVOA CALIBRATION

37	10.451	1258	1269	1276	rVV3	1896135	4498230	5.54%	2.016%
38	10.516	1276	1280	1285	rVB3	353698	609501	0.75%	0.273%
39	10.575	1285	1290	1296	rBV3	998905	1858210	2.29%	0.833%
40	10.839	1329	1335	1343	rBV	1298506	1975397	2.43%	0.885%
41	10.963	1348	1356	1363	rBV2	347855	802385	0.99%	0.360%
42	11.492	1441	1446	1450	rBV2	579485	1010068	1.24%	0.453%
43	11.557	1450	1457	1465	rVB	1895027	3641216	4.49%	1.632%
44	11.728	1478	1486	1495	rBV2	922042	1705896	2.10%	0.764%
45	11.886	1506	1513	1517	rBV2	1066903	1907594	2.35%	0.855%
46	11.928	1517	1520	1525	rVB	803038	1073396	1.32%	0.481%
47	12.028	1533	1537	1539	rBV	363484	510222	0.63%	0.229%
48	12.133	1548	1555	1564	rVB2	1863886	3313509	4.08%	1.485%
49	12.533	1618	1623	1635	rBV4	446777	920632	1.13%	0.413%
50	12.722	1648	1655	1661	rBV4	397794	817133	1.01%	0.366%
51	13.057	1703	1712	1719	rBV2	546185	970255	1.20%	0.435%
52	13.139	1719	1726	1733	rBV	1490445	2208518	2.72%	0.990%
53	13.269	1738	1748	1757	rBV	730008	1598228	1.97%	0.716%
54	13.486	1777	1785	1794	rBV4	445422	1390671	1.71%	0.623%
55	13.575	1795	1800	1805	rBV	1193189	1675915	2.07%	0.751%
56	13.633	1805	1810	1812	rVV2	343956	634087	0.78%	0.284%
57	13.669	1813	1816	1830	rVB6	614582	1911118	2.36%	0.856%
58	13.804	1836	1839	1844	rVB3	413514	649577	0.80%	0.291%
59	13.945	1859	1863	1867	rBV	430405	586524	0.72%	0.263%
60	14.069	1876	1884	1891	rBV4	529387	1258145	1.55%	0.564%
61	14.227	1906	1911	1915	rVB	6269351	8027582	9.89%	3.597%
62	14.316	1922	1926	1931	rVB2	910505	1349187	1.66%	0.605%
63	14.404	1936	1941	1947	rVV5	584903	1055652	1.30%	0.473%
64	14.463	1947	1951	1958	rVV	3423259	4754240	5.86%	2.130%
65	14.527	1958	1962	1970	rVB3	818691	1499495	1.85%	0.672%
66	14.710	1984	1993	2001	rVB7	615039	1774297	2.19%	0.795%
67	14.886	2020	2023	2026	rVV	721748	881643	1.09%	0.395%
68	14.951	2026	2034	2039	rVV	3302716	5566185	6.86%	2.494%
69	14.992	2039	2041	2047	rVB2	352968	491878	0.61%	0.220%
70	15.092	2049	2058	2062	rBV	1349809	2269184	2.80%	1.017%
71	15.163	2066	2070	2071	rVV2	625505	765031	0.94%	0.343%
72	15.186	2071	2074	2081	rVB	1269930	1902438	2.34%	0.852%
73	15.257	2082	2086	2091	rBV4	264041	485508	0.60%	0.218%
74	15.433	2111	2116	2121	rBV2	887143	1458411	1.80%	0.653%
75	15.598	2139	2144	2149	rVB2	522975	959929	1.18%	0.430%
76	15.698	2157	2161	2167	rVB2	487336	817679	1.01%	0.366%

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77	16.074	2221	2225	2227	rBV	1818542	2181741	2.69%	0.978%
78	16.251	2252	2255	2263	rVV5	359496	748242	0.92%	0.335%
79	16.421	2279	2284	2287	rBV5	351884	527664	0.65%	0.236%
80	16.598	2309	2314	2323	rBV5	342442	950944	1.17%	0.426%
81	16.798	2336	2348	2352	rVV	37265514	81134832	100.00%	36.354%
82	16.880	2352	2362	2366	rVV	1341783	2929863	3.61%	1.313%
83	16.933	2366	2371	2382	rVB4	736362	1673071	2.06%	0.750%
84	17.127	2399	2404	2410	rVV	1619363	2859424	3.52%	1.281%
85	17.210	2414	2418	2423	rVV	1330134	1906765	2.35%	0.854%
86	17.257	2423	2426	2433	rVB	965341	1283780	1.58%	0.575%
87	17.321	2433	2437	2441	rBV2	340972	509159	0.63%	0.228%
88	17.421	2449	2454	2462	rVB4	602764	1242109	1.53%	0.557%
89	17.521	2463	2471	2480	rBV7	235932	808954	1.00%	0.362%
90	17.639	2487	2491	2497	rVB	1067562	1464284	1.80%	0.656%
91	17.821	2518	2522	2526	rBV	403177	556051	0.69%	0.249%
92	18.351	2608	2612	2618	rVB	853765	1126122	1.39%	0.505%
93	18.839	2689	2695	2701	rVB5	230476	585879	0.72%	0.263%
94	19.027	2721	2727	2733	rVB4	686858	1621699	2.00%	0.727%
95	19.621	2824	2828	2838	rVB2	443194	699625	0.86%	0.313%
96	20.180	2919	2923	2932	rBV3	870377	1273167	1.57%	0.570%
97	21.221	3096	3100	3105	rVB2	438033	618270	0.76%	0.277%
98	21.539	3149	3154	3161	rVB	1267563	2002061	2.47%	0.897%
99	22.415	3298	3303	3309	rVB2	359014	691023	0.85%	0.310%
100	24.833	3706	3714	3724	rVB	901408	2329724	2.87%	1.044%

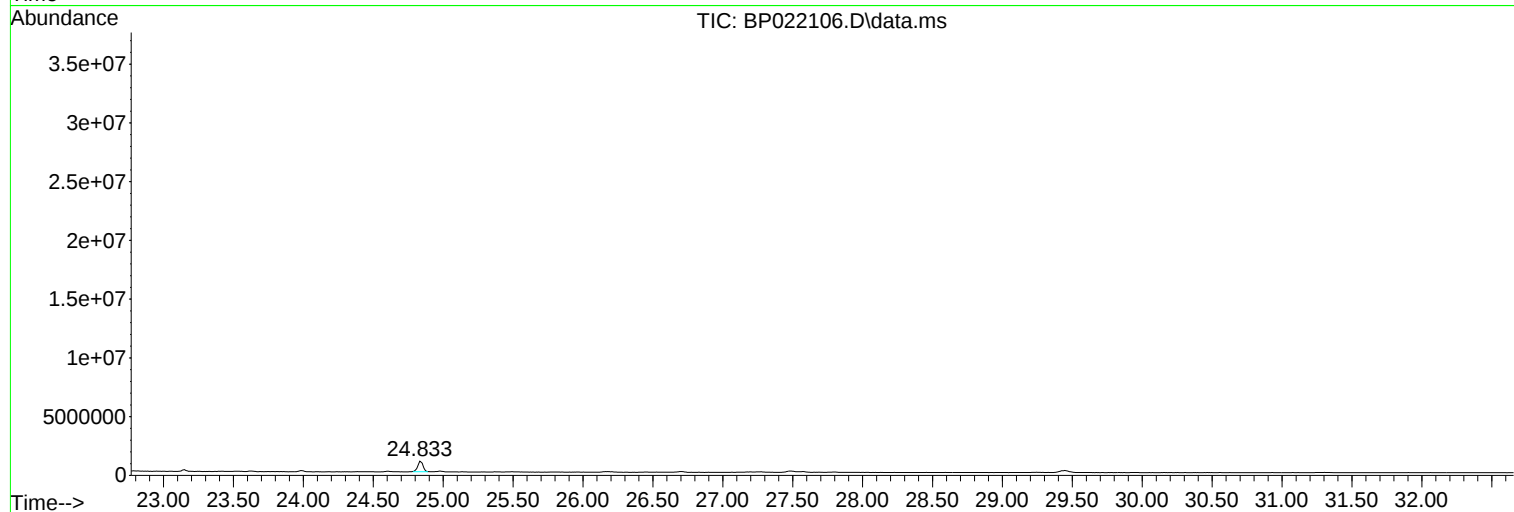
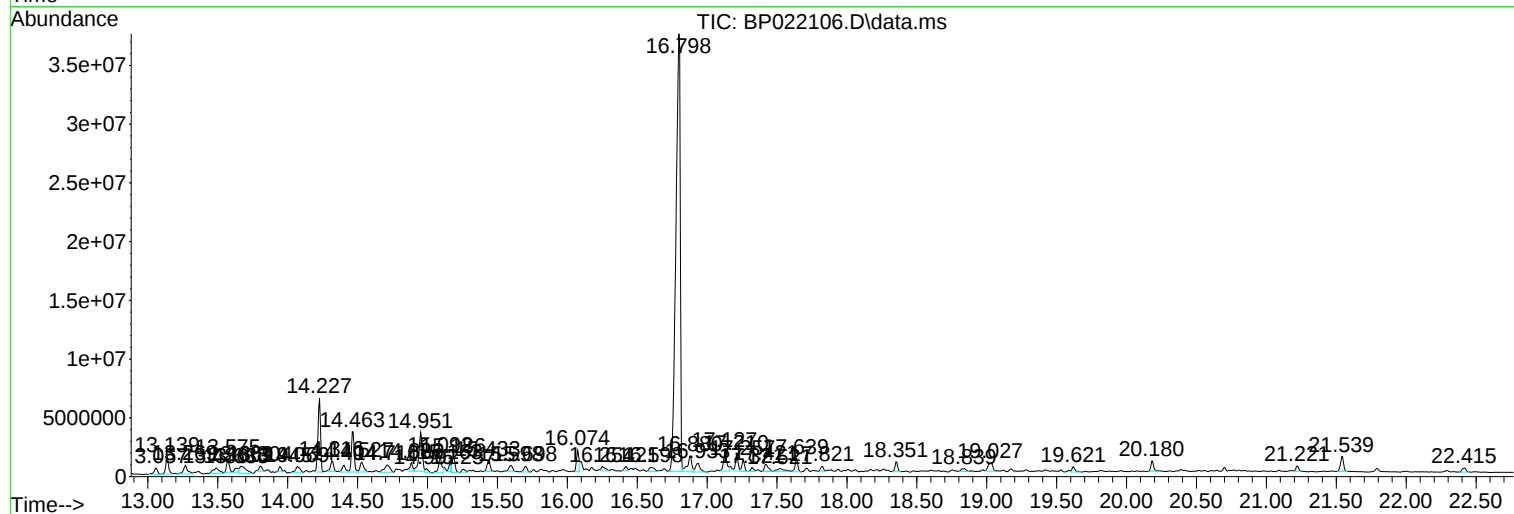
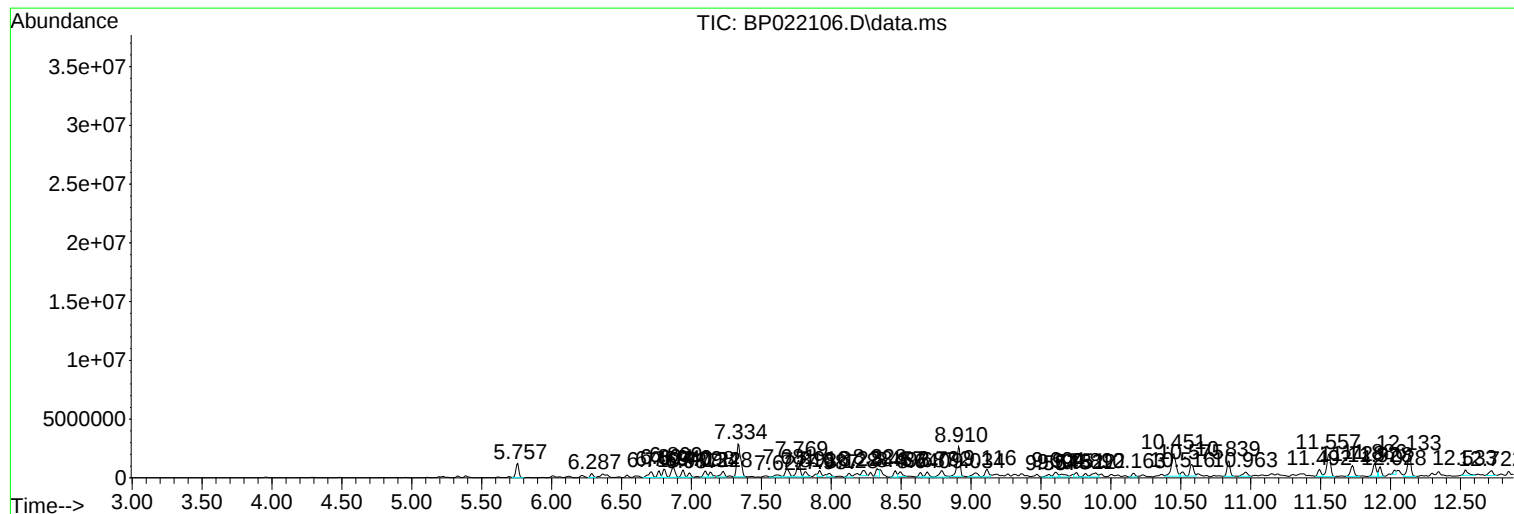
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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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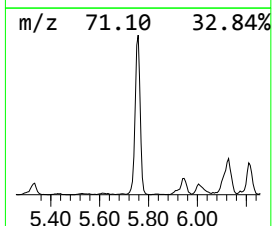
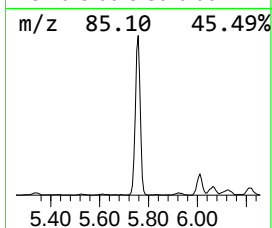
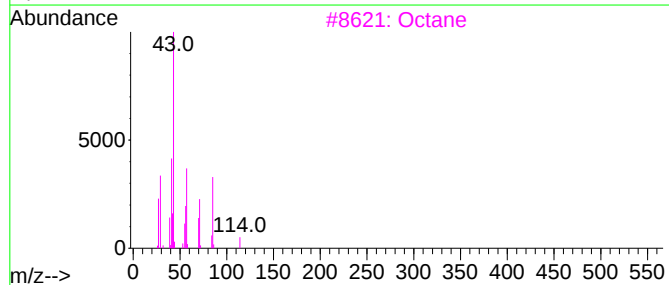
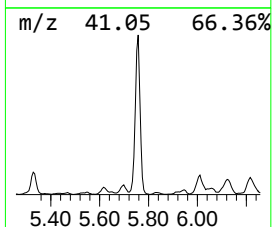
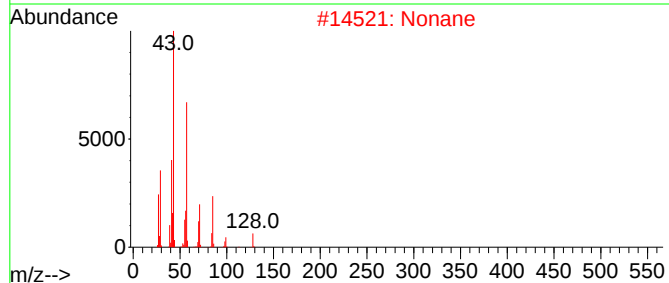
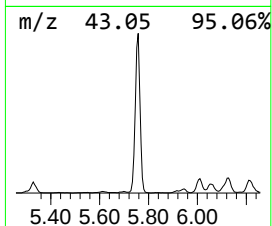
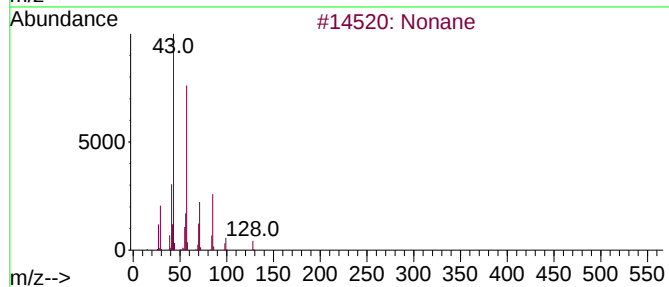
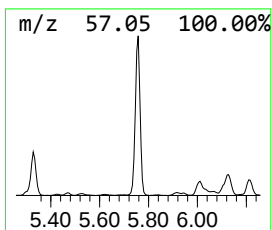
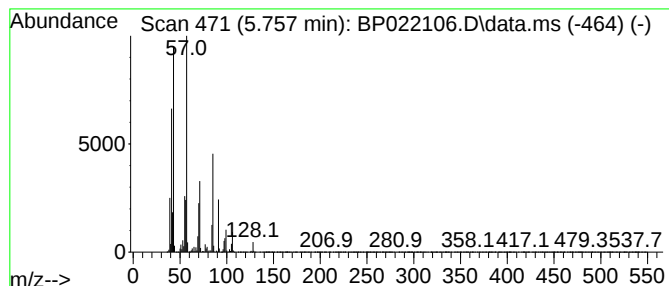
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.757	26.58 ng/ul	1815390	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	93
2			Nonane	128	C9H20	000111-84-2	93
3			Octane	114	C8H18	000111-65-9	59
4			Octane	114	C8H18	000111-65-9	59
5			Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	50



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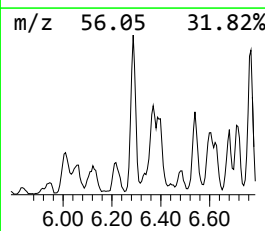
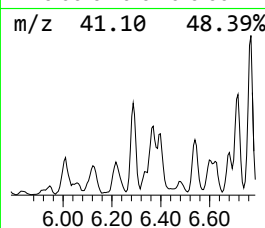
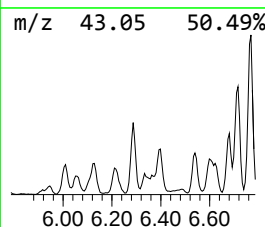
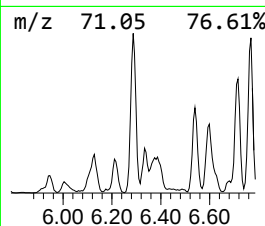
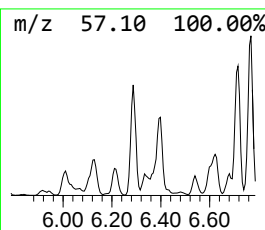
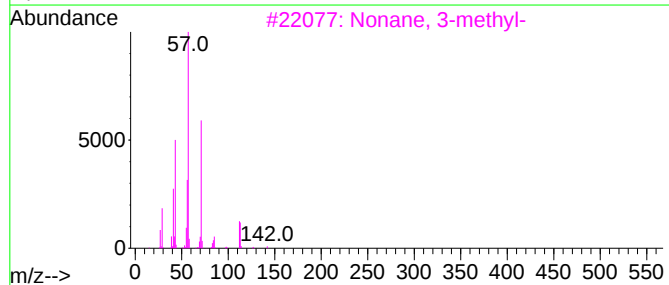
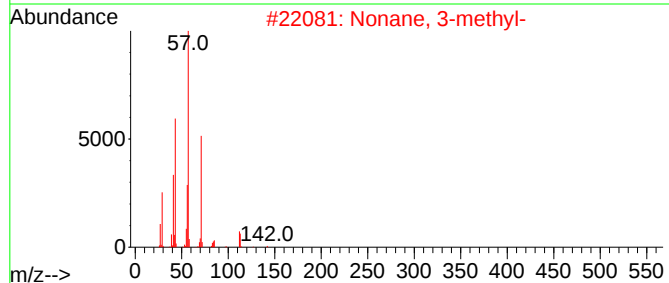
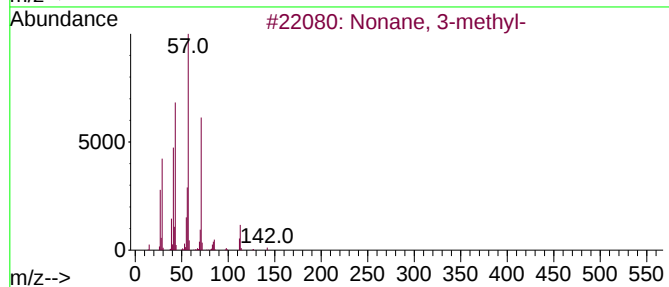
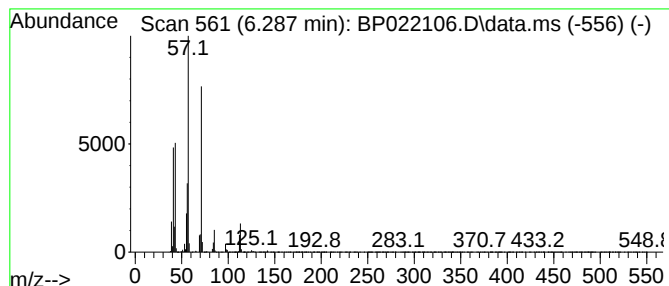
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 57

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.287	7.03 ng/ul	480035	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3-methyl-	142	C10H22	005911-04-6	91
2			Nonane, 3-methyl-	142	C10H22	005911-04-6	90
3			Nonane, 3-methyl-	142	C10H22	005911-04-6	90
4			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	90
5			Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	80



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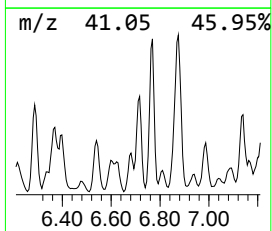
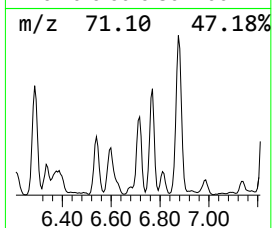
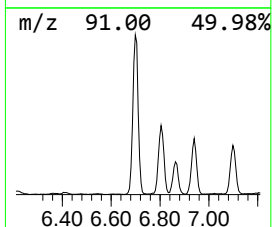
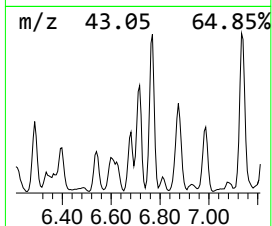
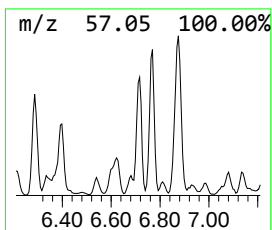
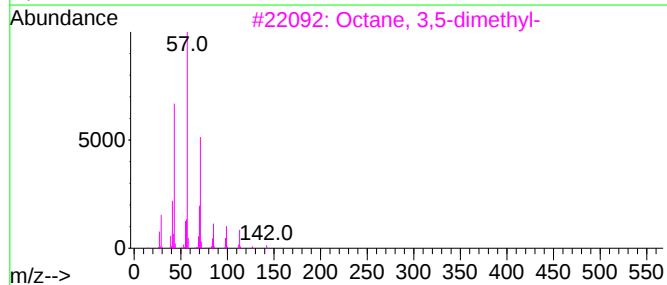
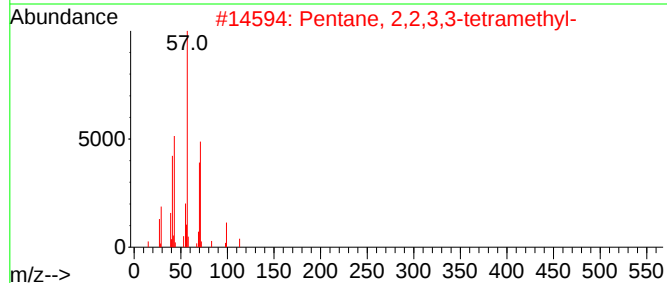
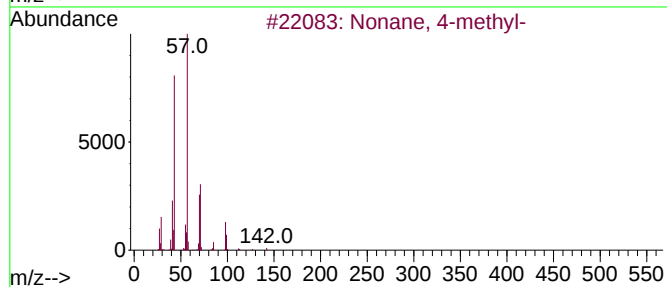
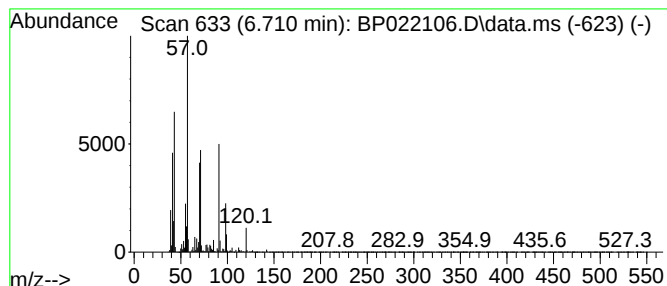
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.710	16.69 ng/ul	1139930	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 4-methyl-	142	C10H22	017301-94-9	70
2			Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	53
3			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	49
4			1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	49
5			Heptane, 4-methyl-	114	C8H18	000589-53-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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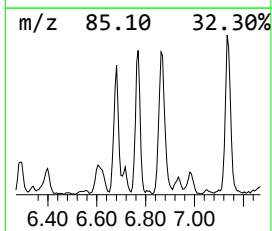
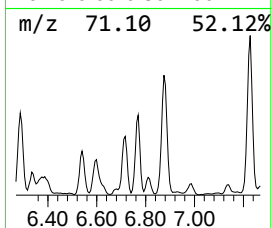
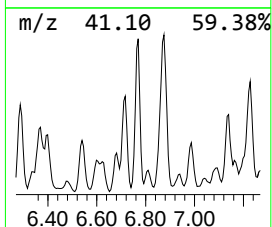
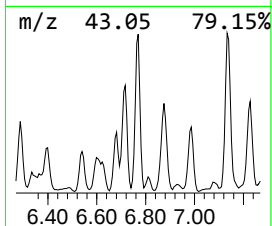
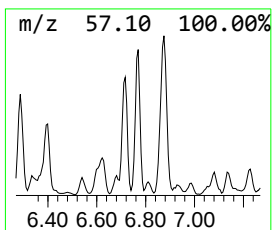
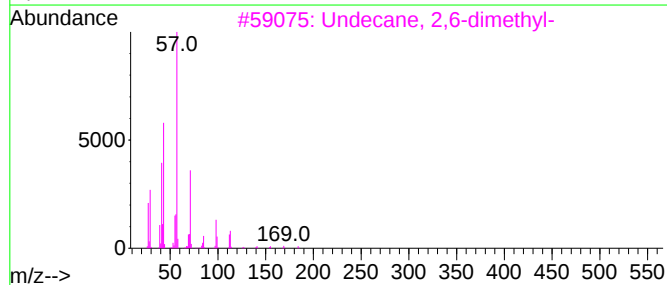
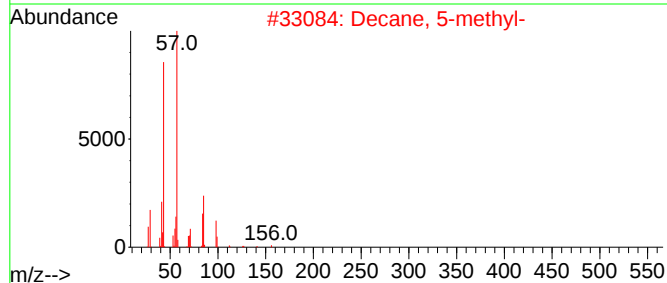
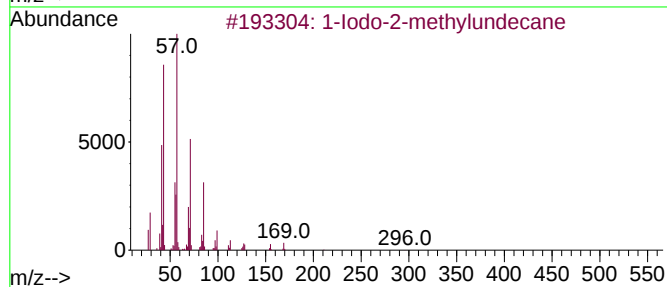
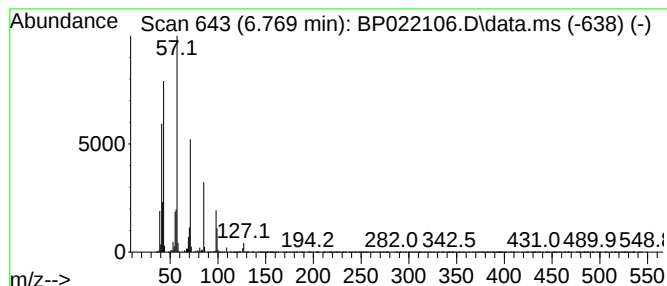
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1-Iodo-2-methylundecane Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.769	11.71 ng/ul	800150	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	64
2			Decane, 5-methyl-	156	C11H24	013151-35-4	53
3			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	53
4			Decane, 2,4-dimethyl-	170	C12H26	002801-84-5	50
5			Undecane, 5-methyl-	170	C12H26	001632-70-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
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Misc :
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Instrument :
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ClientSampleId :
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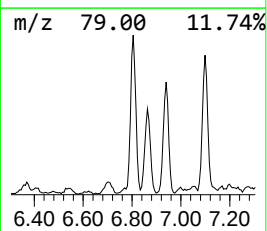
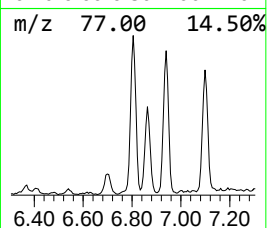
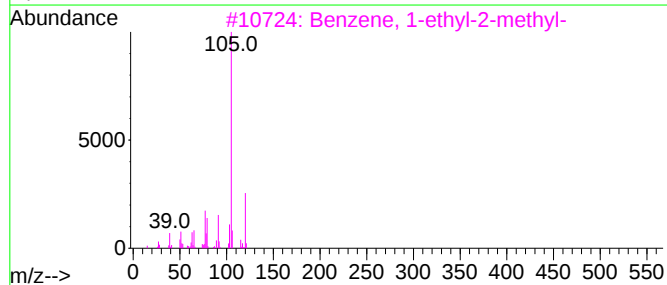
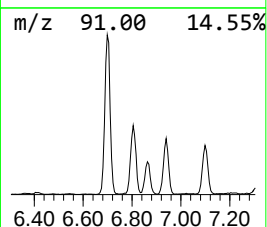
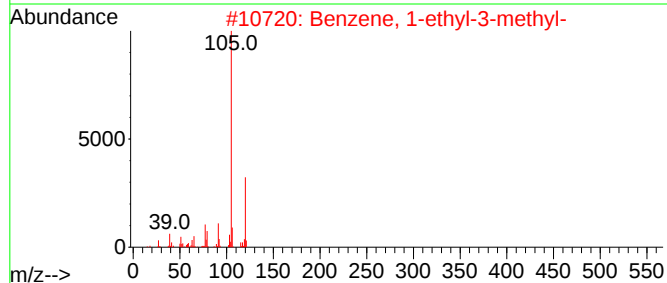
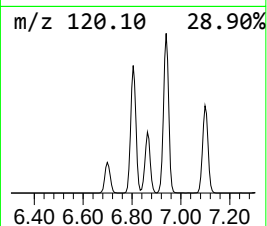
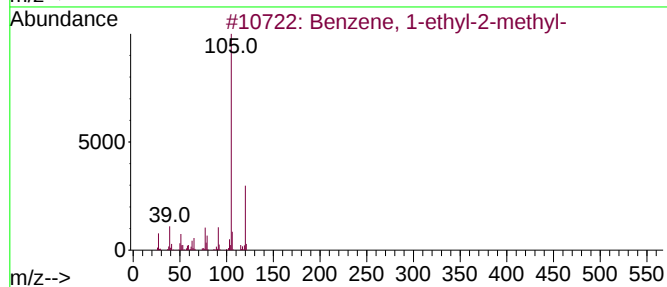
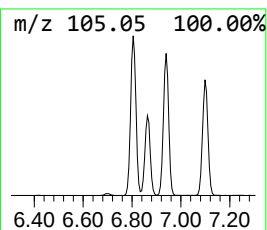
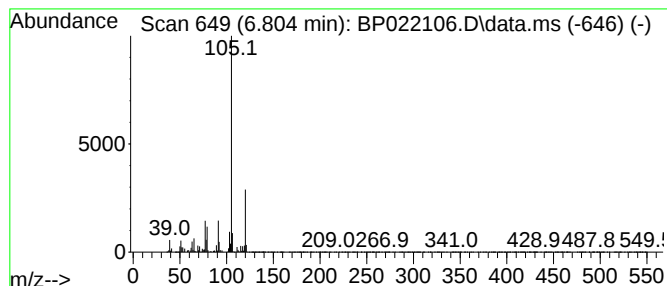
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.804	14.97 ng/ul	1022430	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
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Misc :
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Instrument :
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ClientSampleId :
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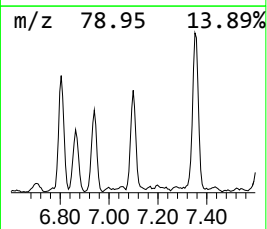
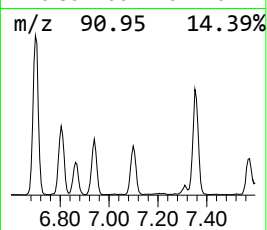
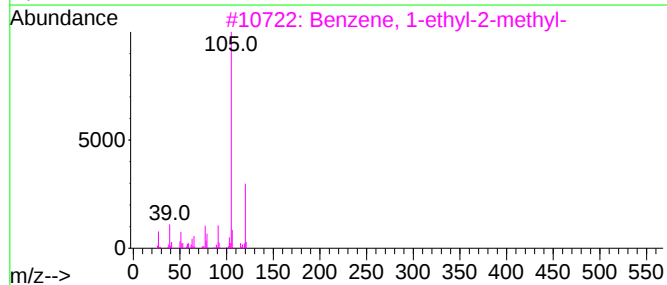
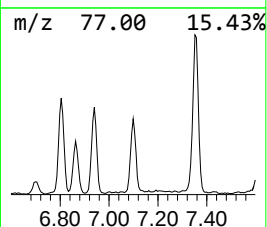
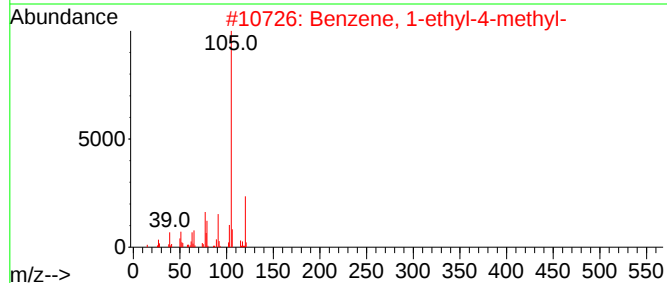
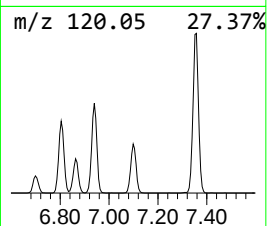
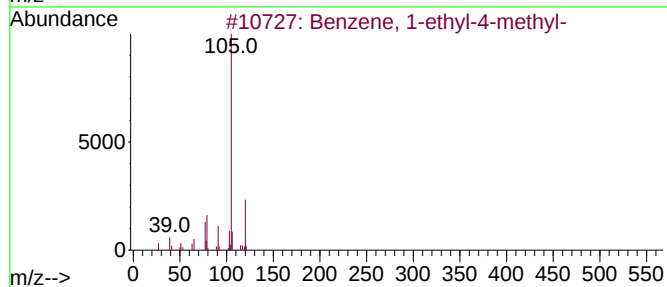
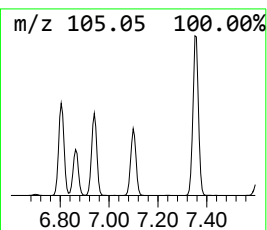
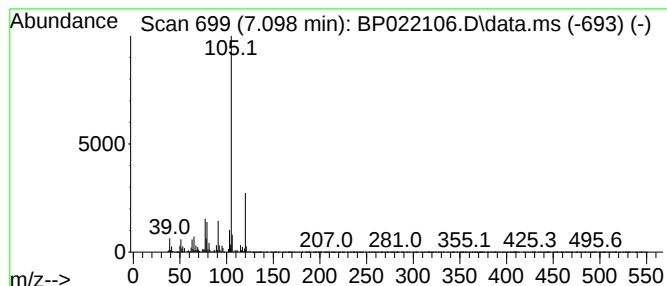
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1-ethyl-4-methyl- Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.098	12.06 ng/ul	824139	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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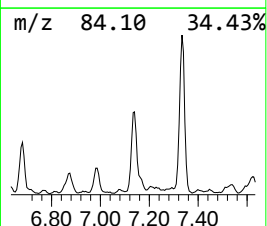
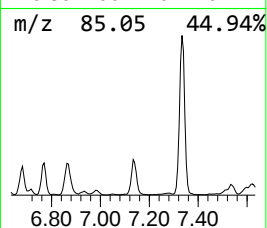
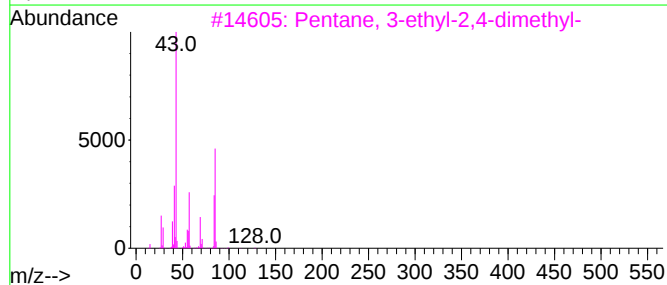
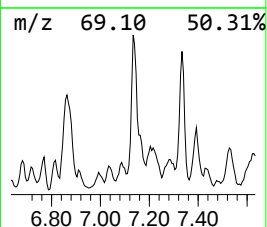
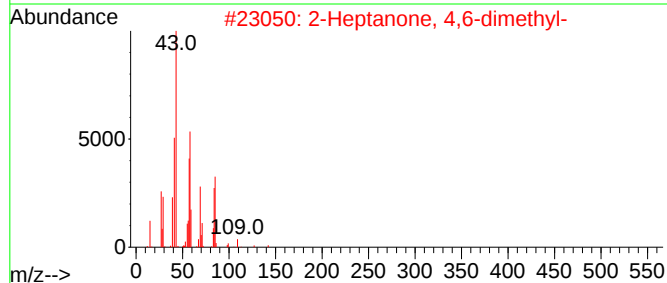
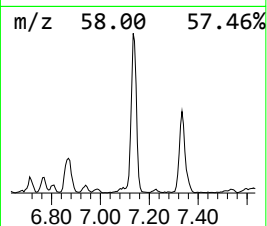
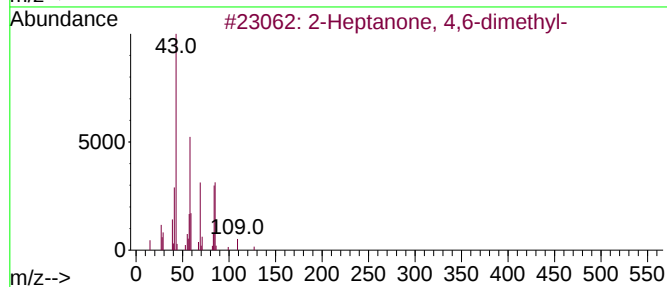
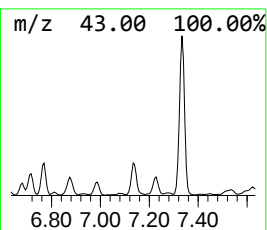
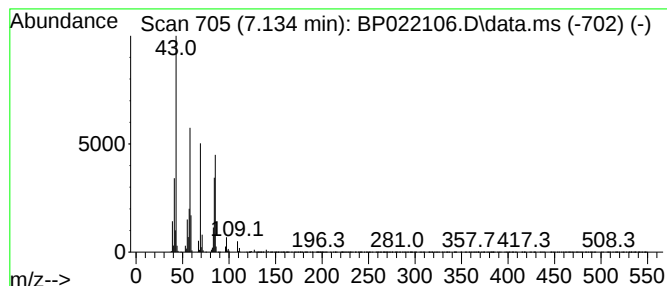
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 2-Heptanone, 4,6-dimethyl- Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.134	9.14 ng/ul	624488	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	80		
2	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	56		
3	Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	49		
4	Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	43		
5	Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
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Misc :
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Instrument :
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ClientSampleId :
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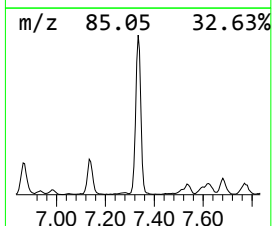
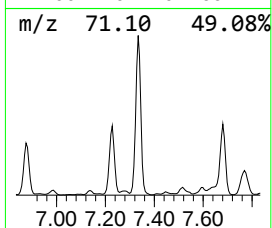
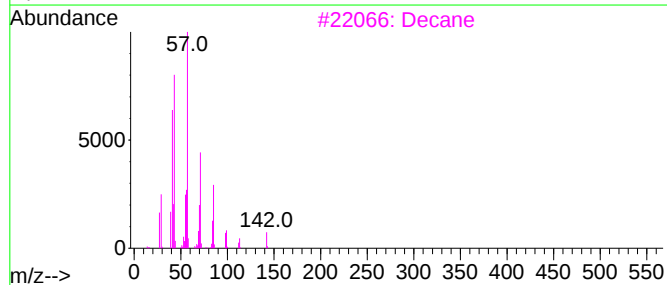
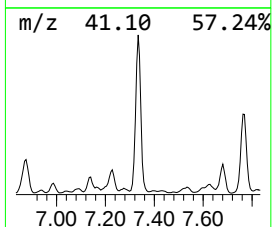
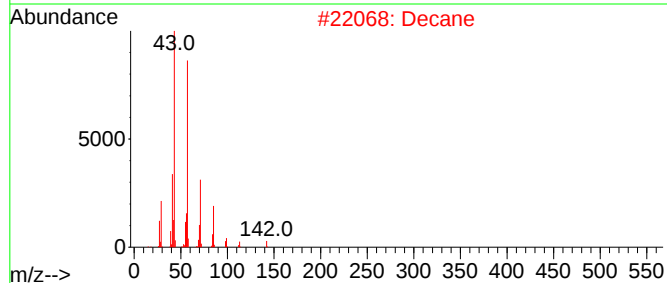
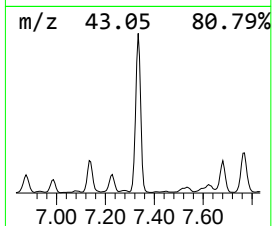
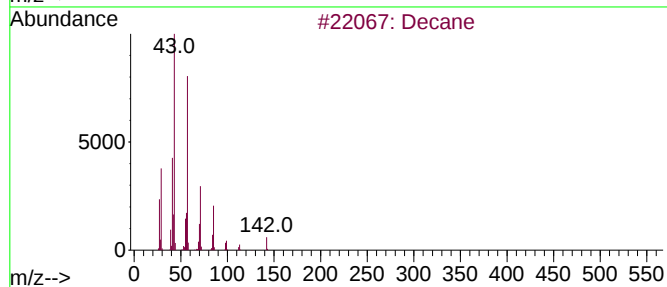
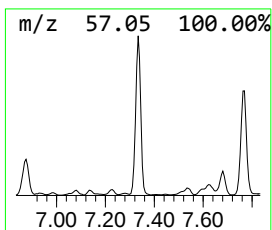
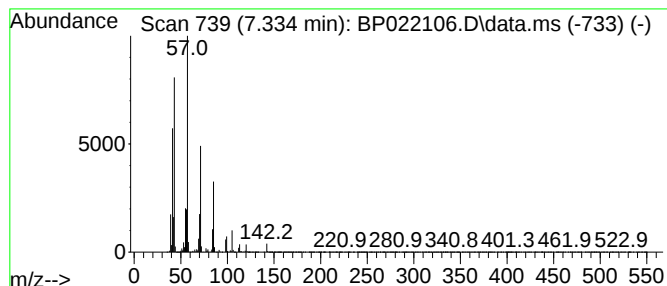
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.334	80.36 ng/ul	5489660	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	93
2			Decane	142	C10H22	000124-18-5	93
3			Decane	142	C10H22	000124-18-5	90
4			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	64
5			Hexadecane	226	C16H34	000544-76-3	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
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Misc :
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Instrument :
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ClientSampleId :
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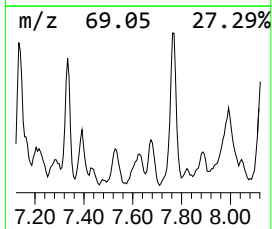
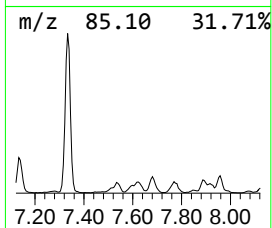
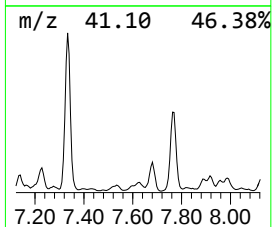
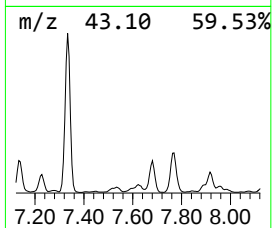
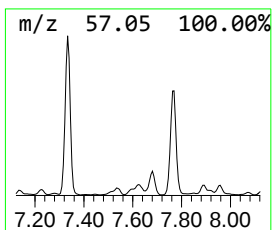
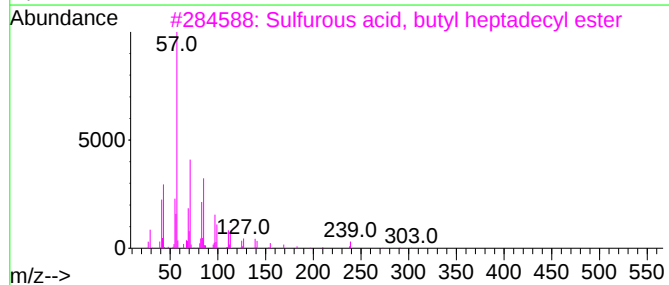
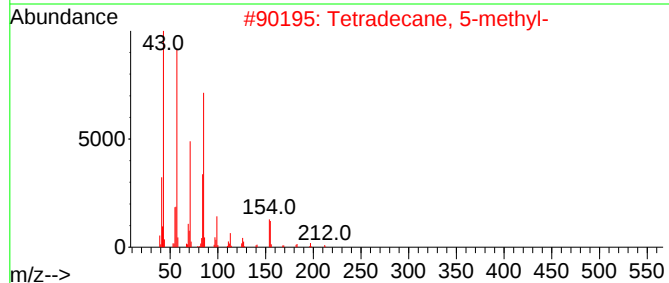
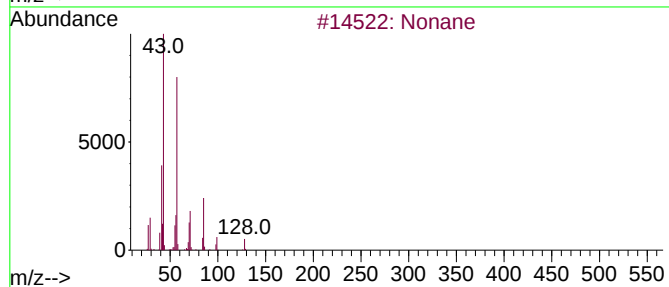
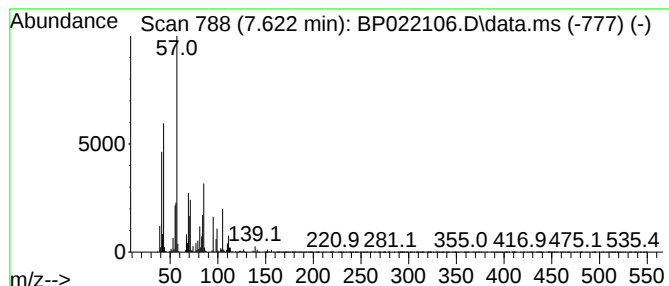
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.622	8.42 ng/ul	575266	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	50
2			Tetradecane, 5-methyl-	212	C15H32	025117-32-2	43
3			Sulfurous acid, butyl heptadecyl...	376	C21H44O3S	1000309-18-4	43
4			Heptane, 2,2,3,3,5,6,6-heptamethyl-	198	C14H30	007225-67-4	43
5			Octane	114	C8H18	000111-65-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
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Misc :
ALS Vial : 12 Sample Multiplier: 1

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ClientSampleId :
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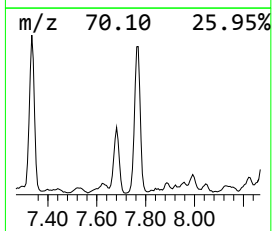
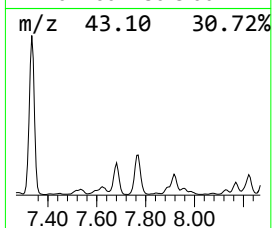
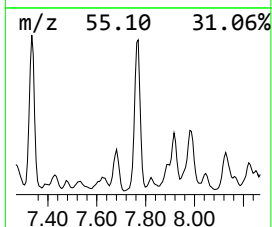
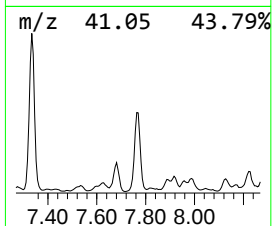
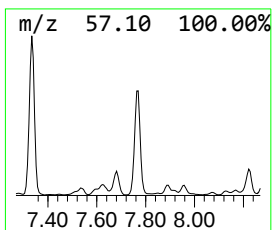
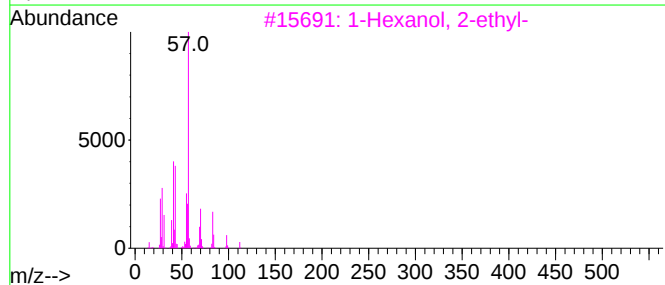
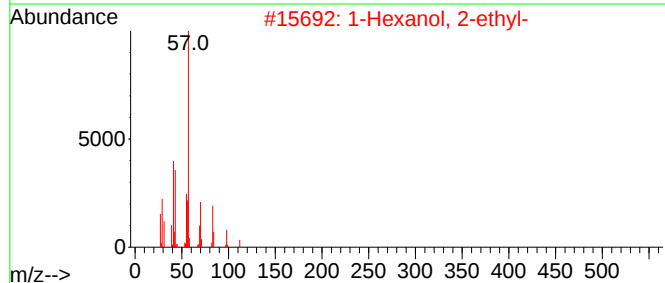
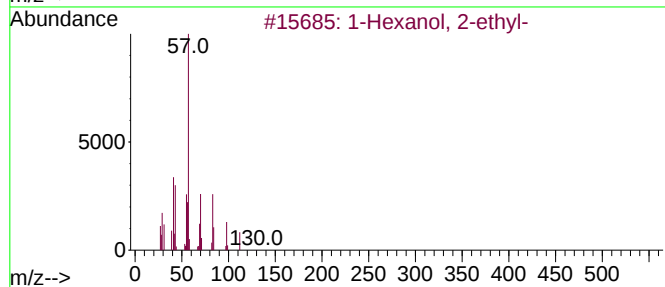
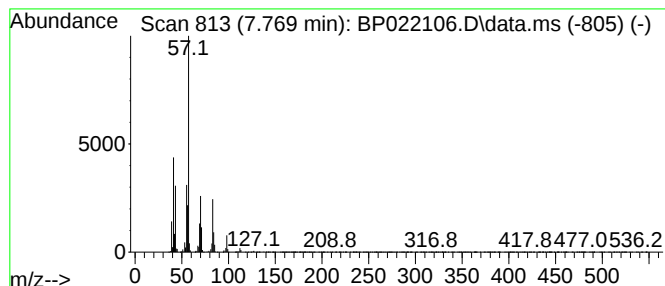
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 1-Hexanol, 2-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.769	34.64 ng/ul	2365960	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
5	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC29DL

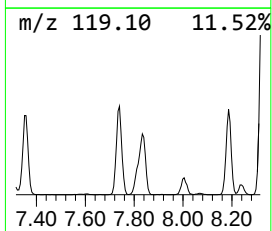
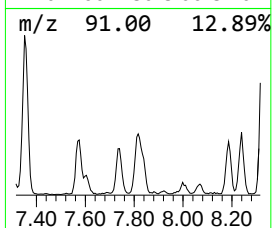
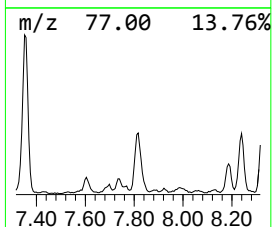
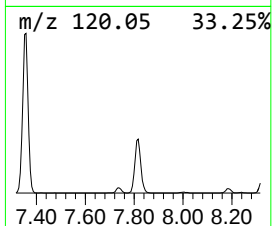
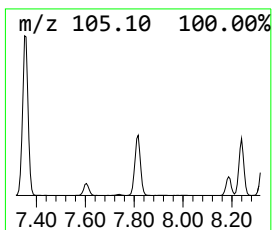
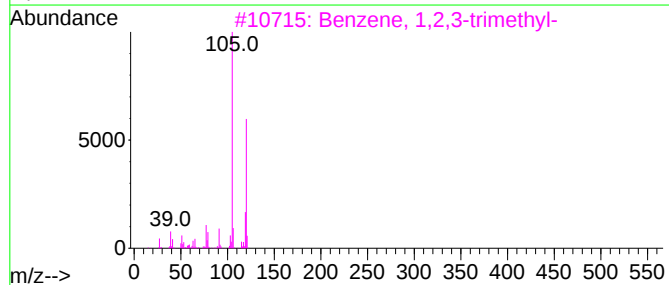
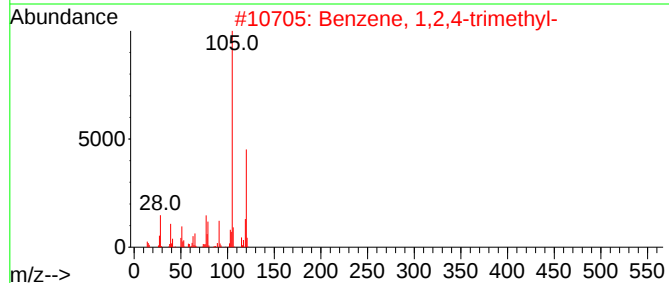
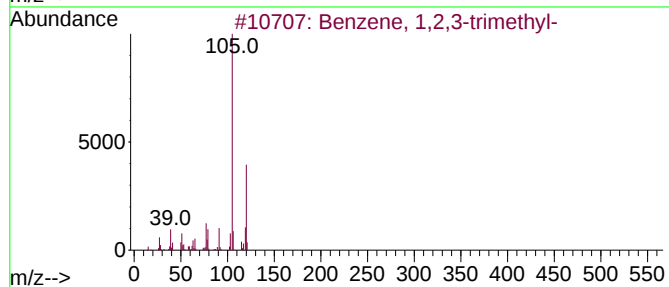
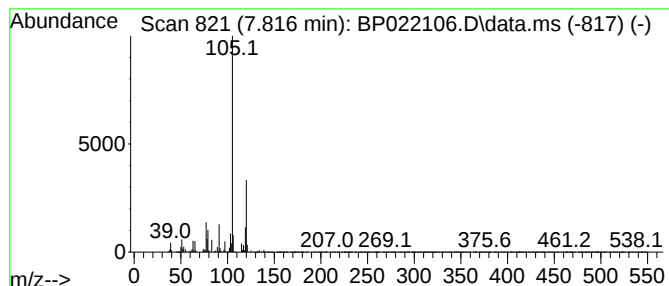
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Benzene, 1,2,3-trimethyl- Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.816	11.32 ng/ul	773555	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
2			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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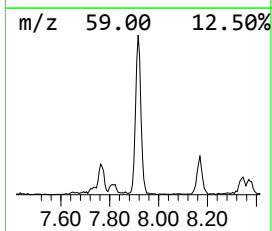
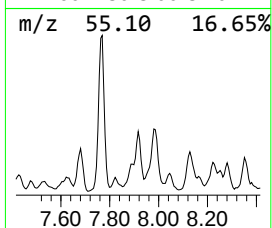
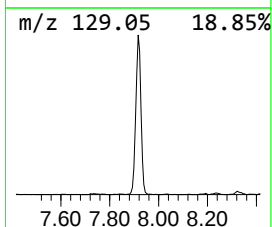
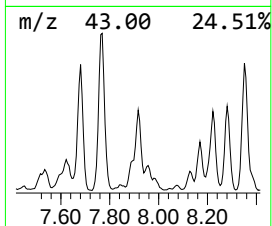
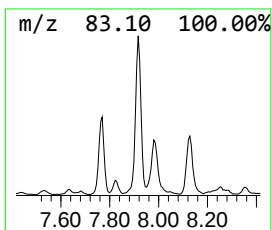
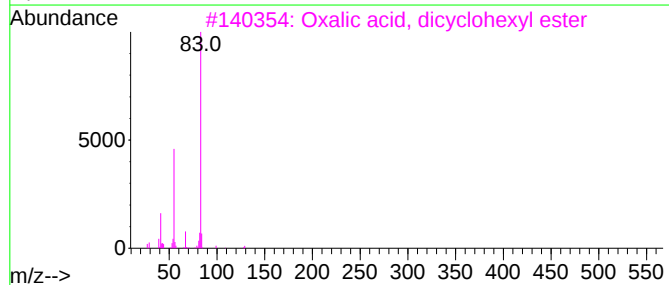
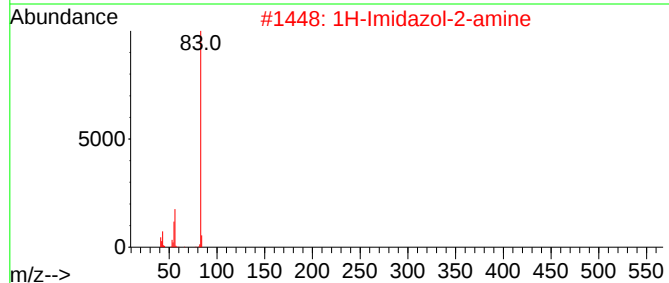
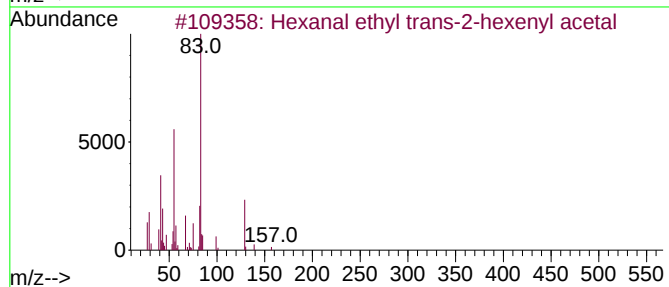
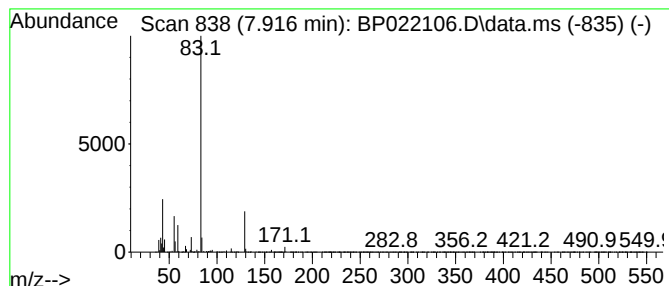
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Hexanal ethyl trans-2-hexen... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.916	10.85 ng/ul	741001	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanal ethyl trans-2-hexenyl ac...	228	C14H28O2	1000431-42-6	50
2			1H-Imidazol-2-amine	83	C3H5N3	007720-39-0	42
3			Oxalic acid, dicyclohexyl ester	254	C14H22O4	1000309-30-9	38
4			N-Nitro-1H-1,2,4-triazol-3-amine	129	C2H3N5O2	034815-01-5	38
5			Oxalic acid, cyclohexyl octyl ester	284	C16H28O4	1000309-31-0	36



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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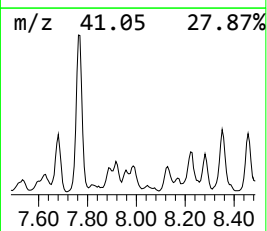
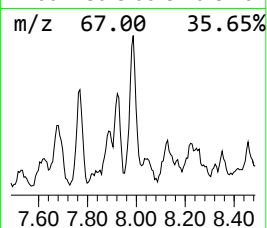
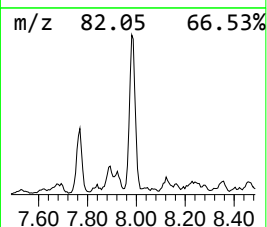
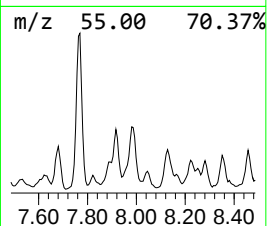
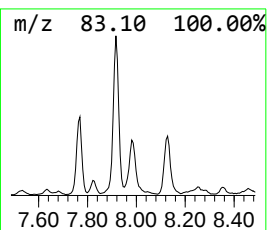
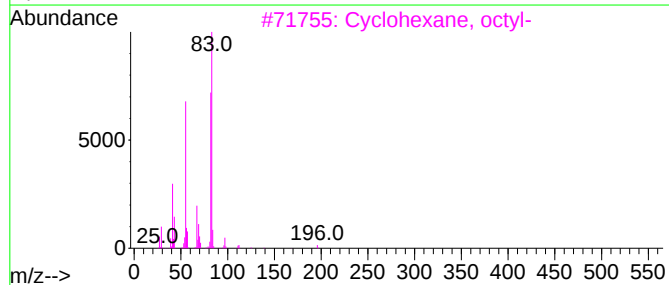
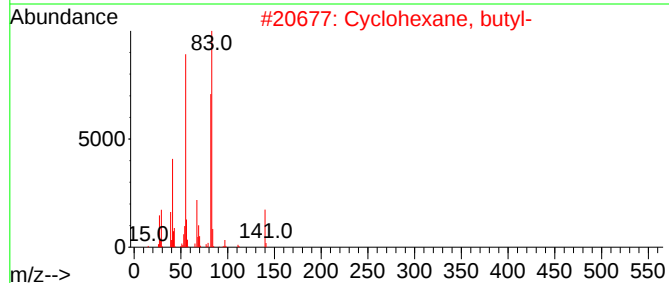
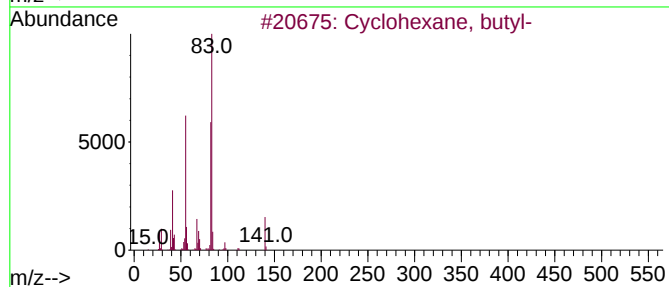
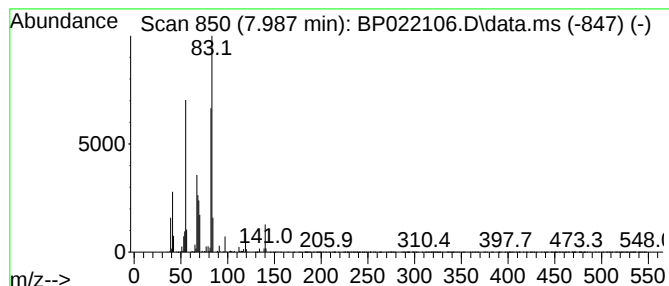
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 (DEL) Alkane: Cyclic7.987 Concentration Rank 56

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.987	7.06 ng/ul	481988	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, butyl-	140	C10H20	001678-93-9	72
2			Cyclohexane, butyl-	140	C10H20	001678-93-9	72
3			Cyclohexane, octyl-	196	C14H28	001795-15-9	64
4			Cyclohexane, decyl-	224	C16H32	001795-16-0	64
5			Cyclohexane, butyl-	140	C10H20	001678-93-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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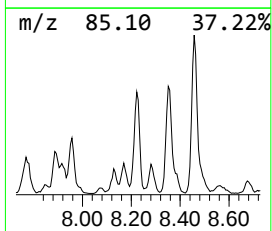
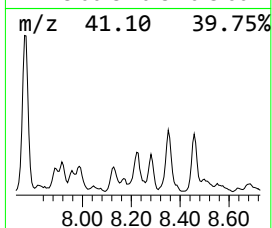
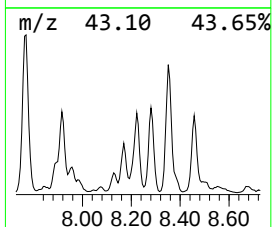
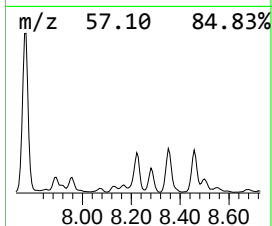
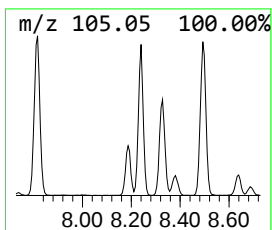
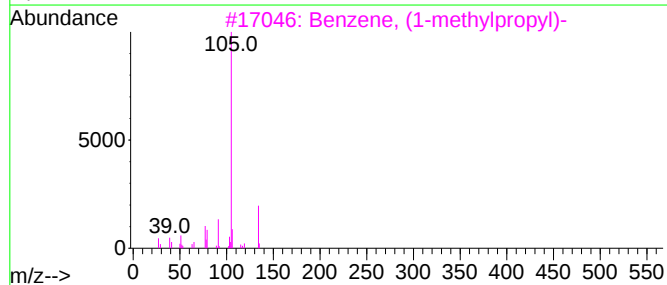
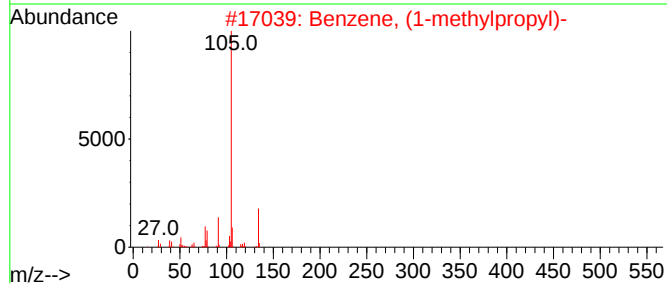
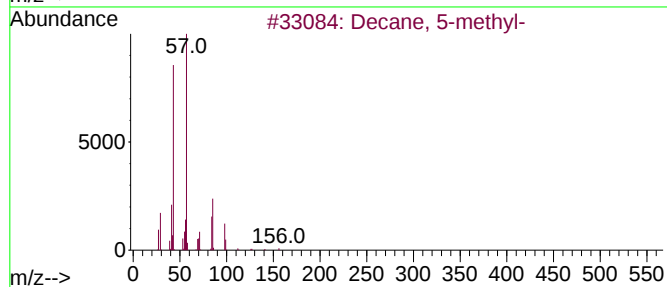
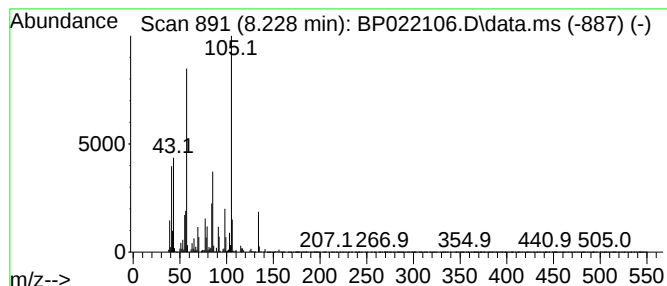
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.228	12.38 ng/ul	845738	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 5-methyl-	156	C11H24	013151-35-4	49
2			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	35
3			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	35
4			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	35
5			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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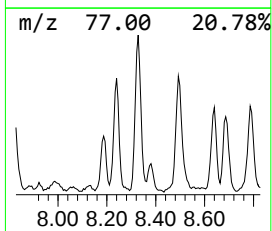
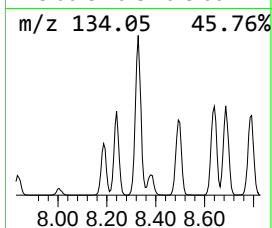
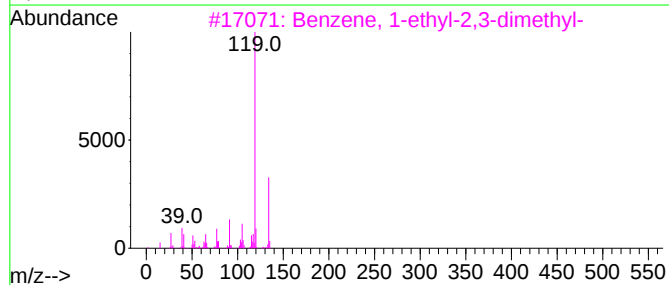
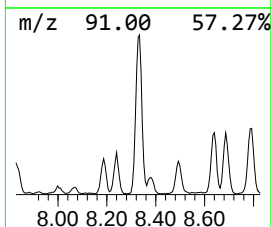
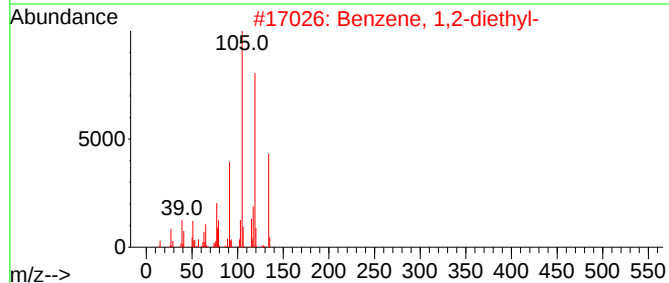
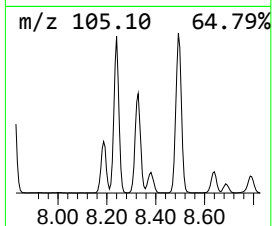
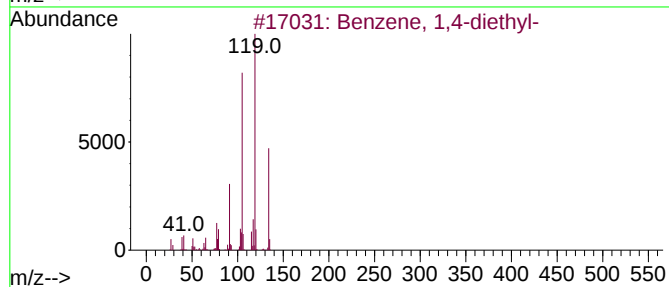
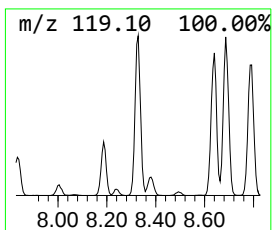
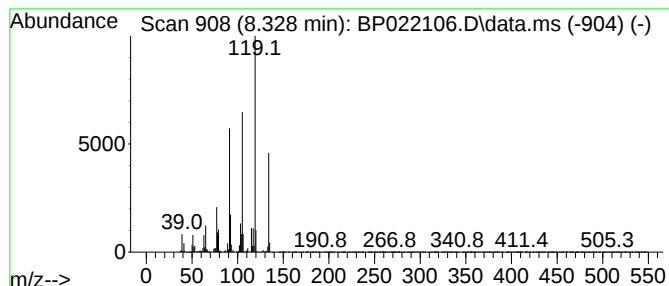
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Benzene, 1,4-diethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.328	16.76 ng/ul	1144660	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	94
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5			o-Cymene	134	C10H14	000527-84-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
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Misc :
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Instrument :
BNA_P
ClientSampleId :
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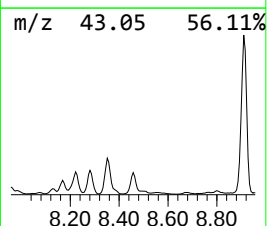
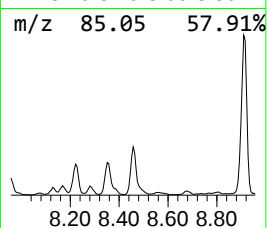
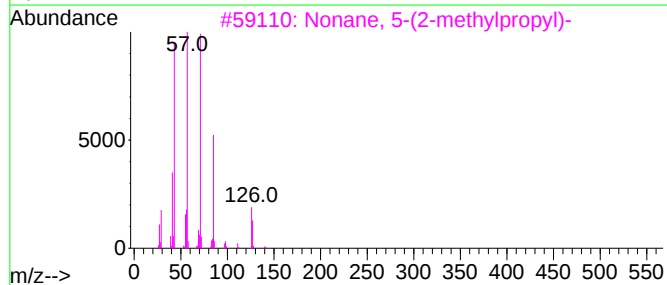
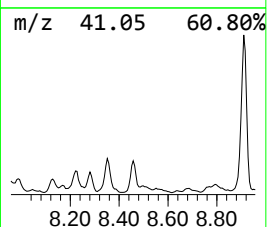
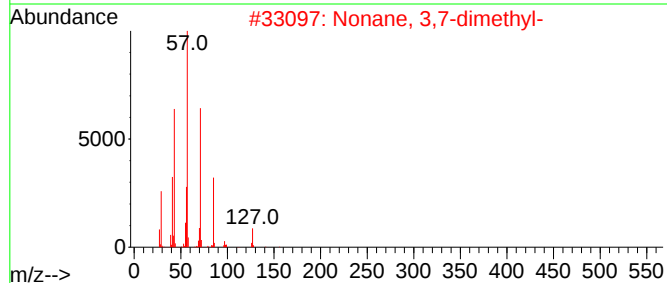
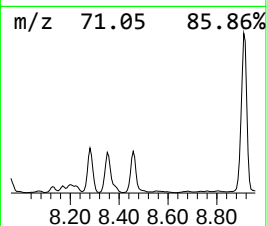
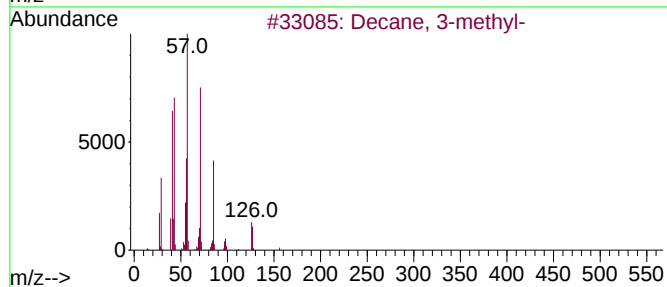
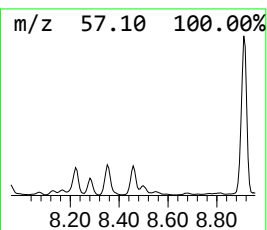
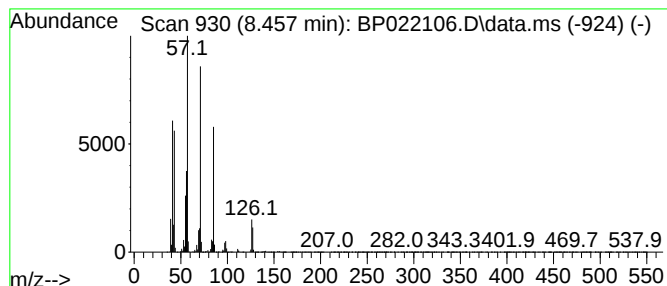
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.457	12.28 ng/ul	838997	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3-methyl-	156	C11H24	013151-34-3	90
2			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	83
3			Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	78
4			Undecane, 2,8-dimethyl-	184	C13H28	017301-25-6	64
5			Nonane, 1-iodo-	254	C9H19I	004282-42-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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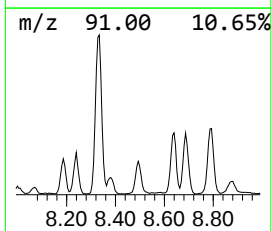
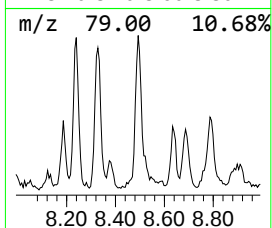
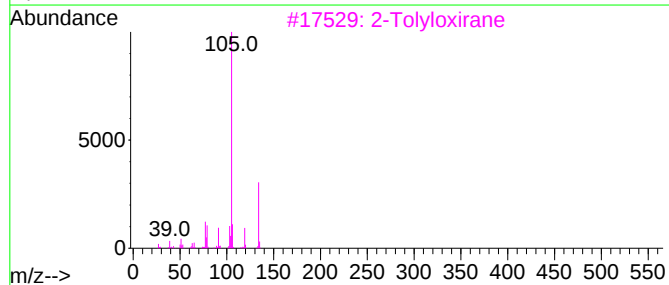
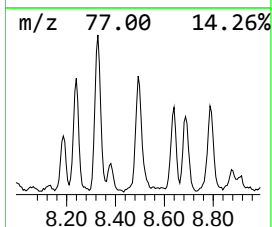
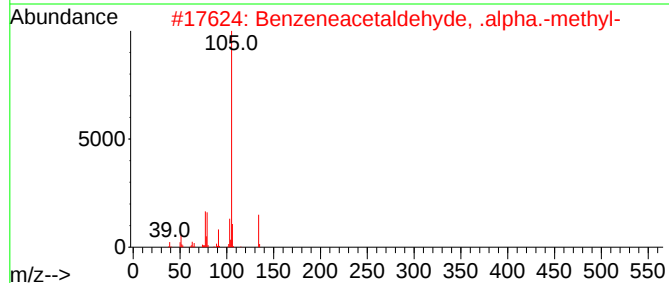
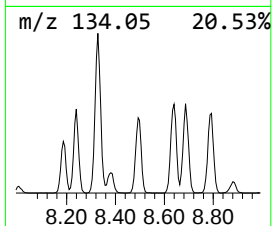
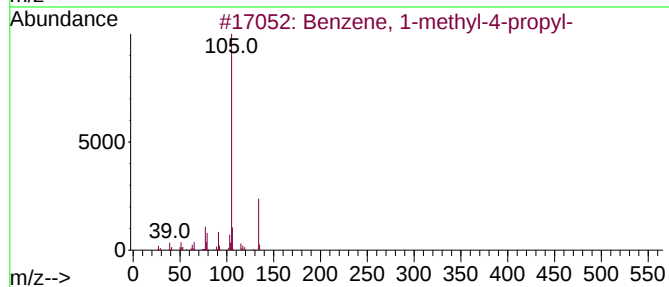
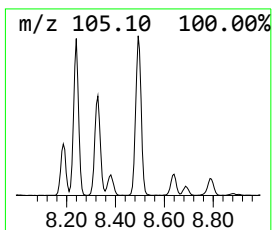
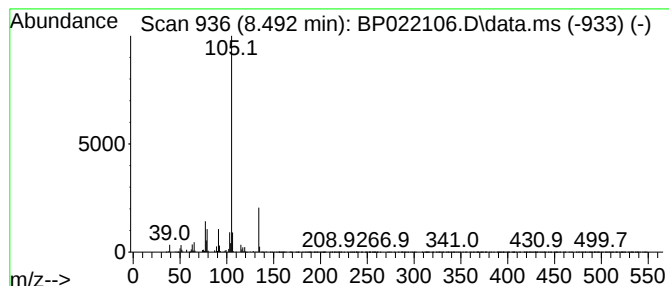
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Benzene, 1-methyl-4-propyl- Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.492	10.68 ng/ul	729664	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	94
2			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	91
3			2-Tolyloxirane	134	C9H10O	002783-26-8	91
4			Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	91
5			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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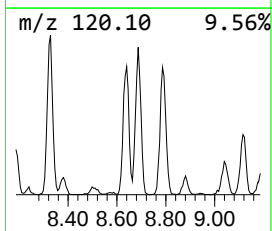
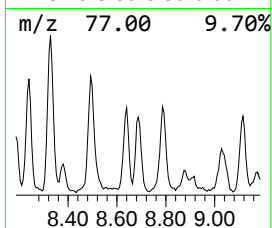
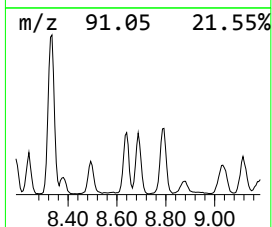
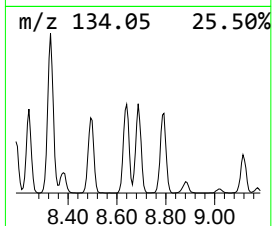
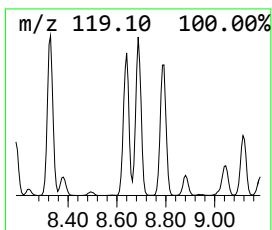
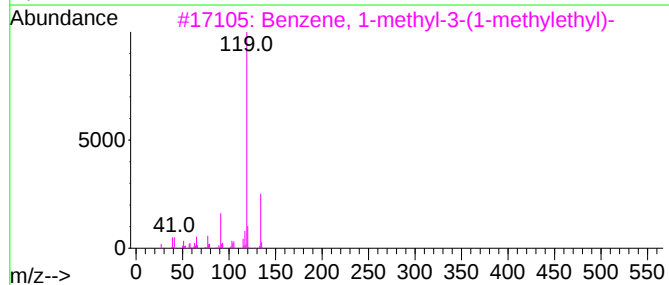
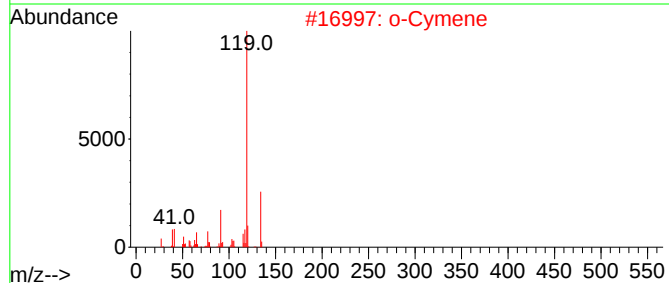
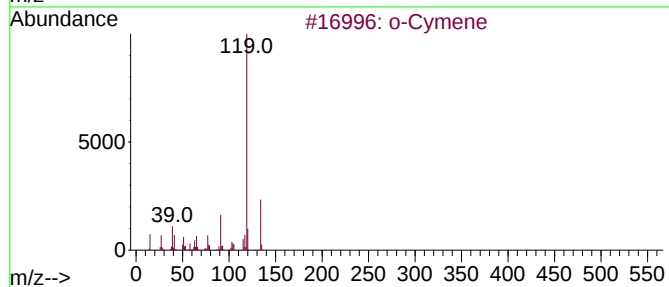
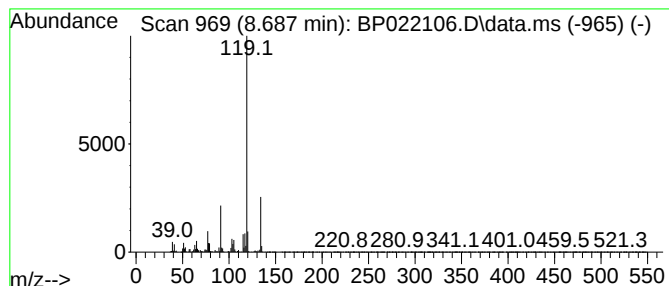
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 o-Cymene Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.687	10.39 ng/ul	709986	1,4-Dichlorobenzene-d4	7.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	97
2			o-Cymene	134	C10H14	000527-84-4	97
3			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	97
4			p-Cymene	134	C10H14	000099-87-6	97
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
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Misc :
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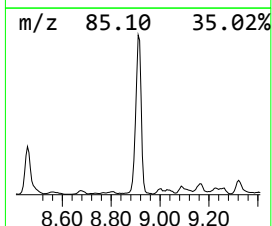
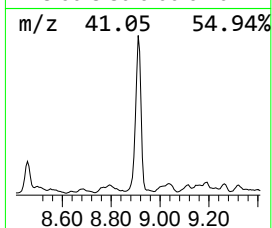
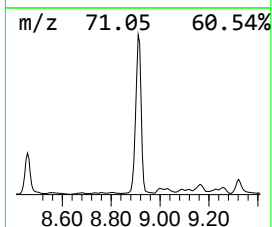
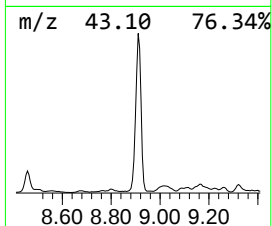
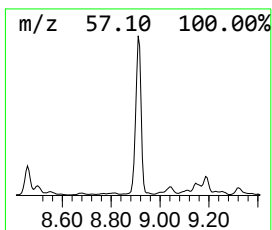
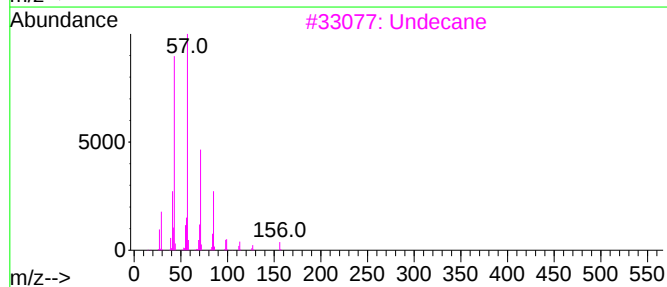
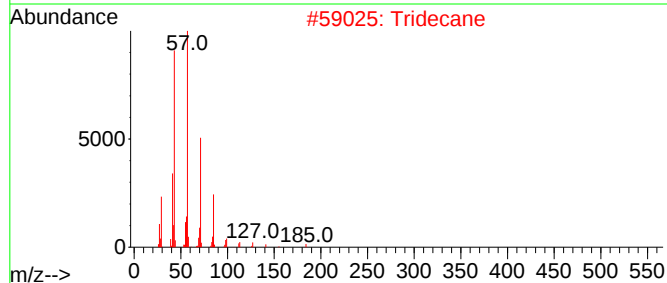
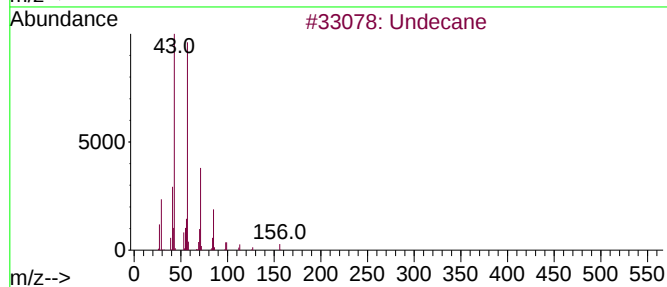
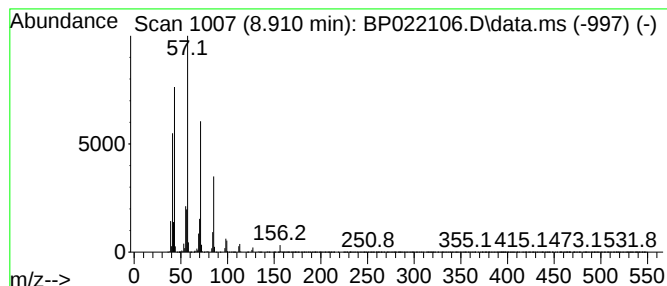
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.910	59.92 ng/ul	4092970	1,4-Dichlorobenzene-d4	7.698

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Undecane	156	C11H24	001120-21-4	90
2		Tridecane	184	C13H28	000629-50-5	90
3		Undecane	156	C11H24	001120-21-4	83
4		Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	83
5		Hexadecane	226	C16H34	000544-76-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
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Sample : P3910-10DL 10X
Misc :
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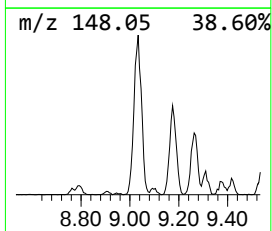
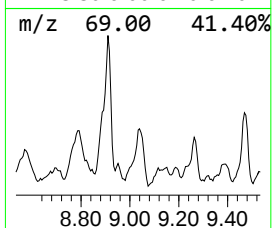
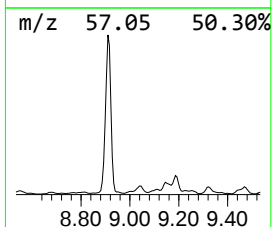
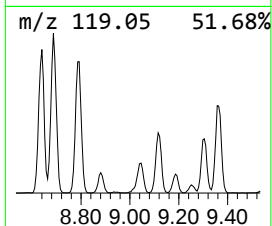
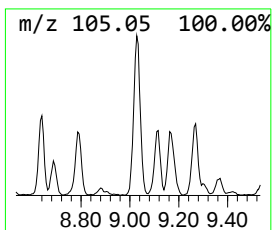
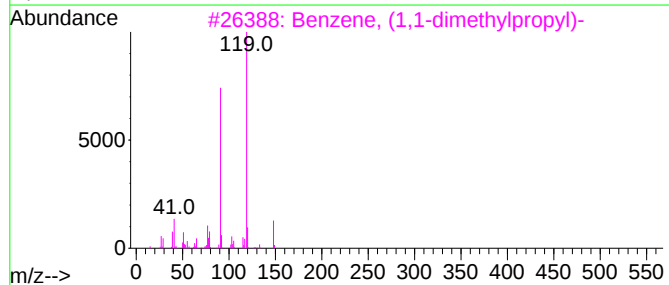
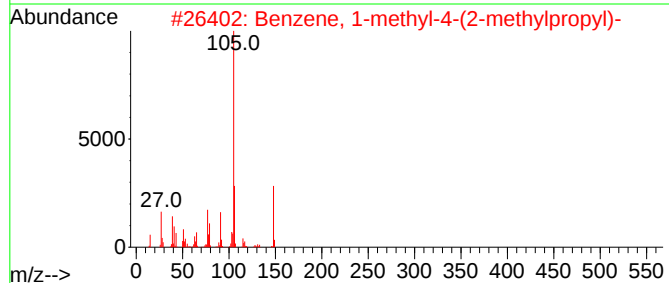
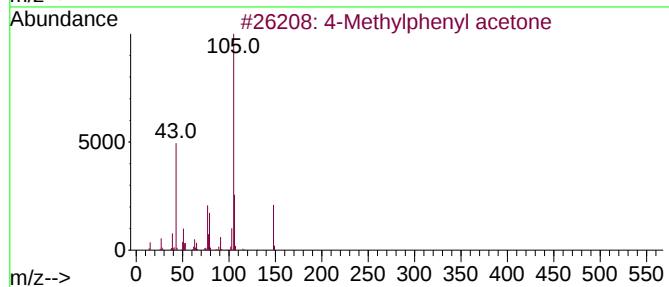
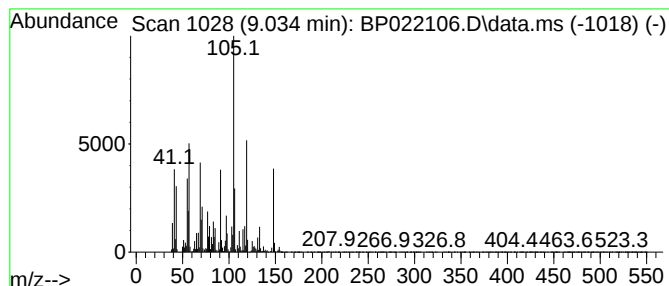
Instrument :
BNA_P
ClientSampleId :
DDC29DL

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 23 unknown-01 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.034	13.47 ng/ul	920244	1,4-Dichlorobenzene-d4	7.698	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methylphenyl acetone	148	C10H12O	002096-86-8	45
2	Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	45
3	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	41
4	2-Butanone, 4-phenyl-	148	C10H12O	002550-26-7	41
5	2-Ethyl[1,3]dithiane	148	C6H12S2	006007-23-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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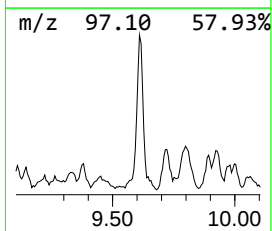
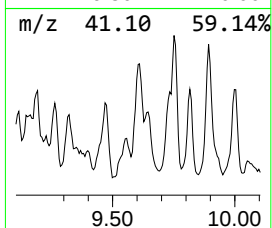
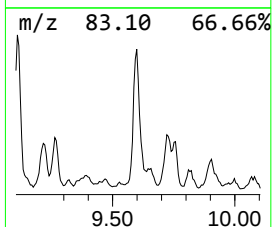
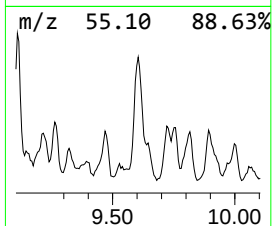
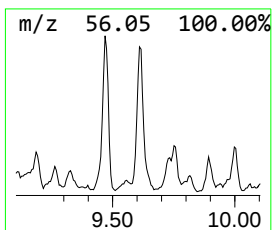
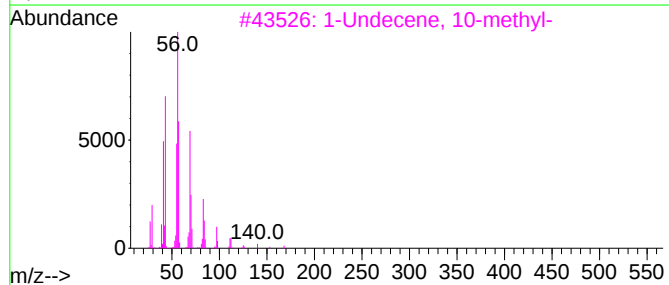
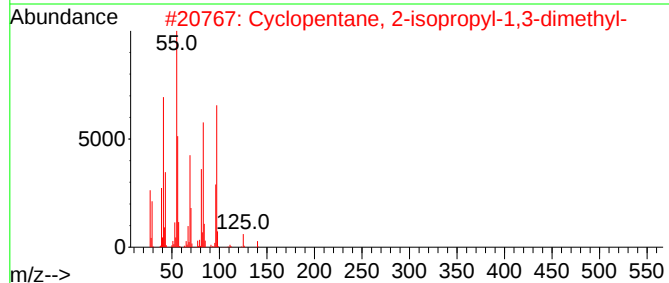
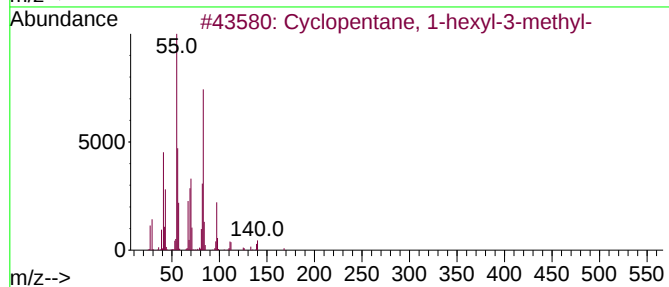
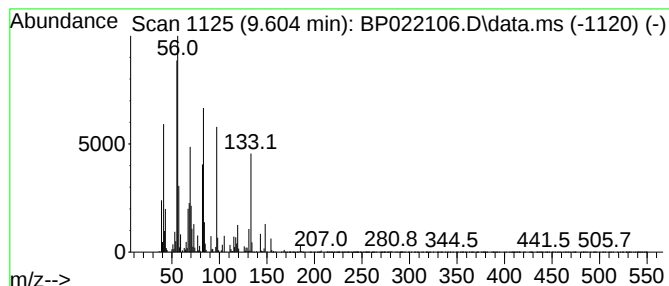
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 25 (DEL) Alkane: Cyclic9.604 Concentration Rank 69

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.604	3.70 ng/ul	832860	Naphthalene-d8	10.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1-hexyl-3-methyl-	168	C12H24	061142-68-5	38
2			Cyclopentane, 2-isopropyl-1,3-di...	140	C10H20	032281-85-9	35
3			1-Undecene, 10-methyl-	168	C12H24	022370-55-4	30
4			2-Decene, 9-methyl-, (Z)-	154	C11H22	074630-24-3	25
5			Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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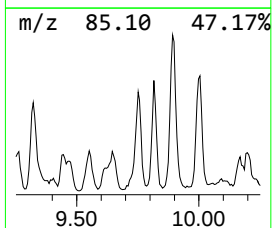
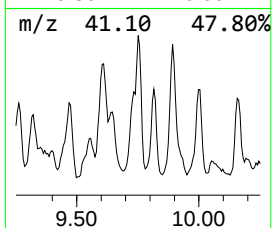
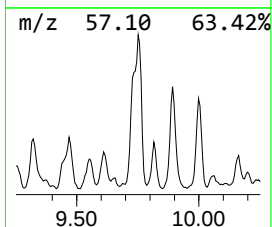
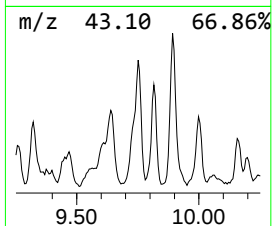
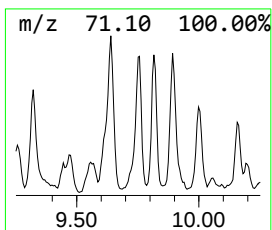
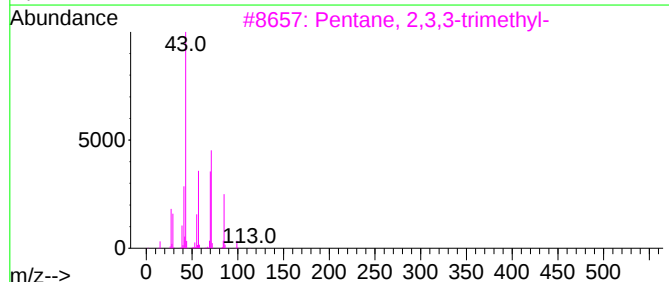
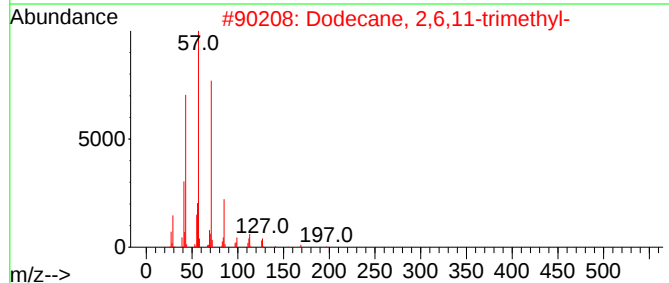
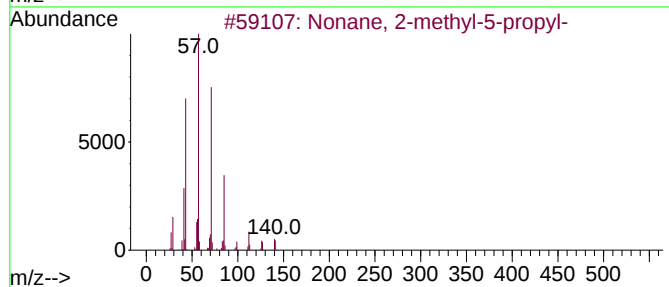
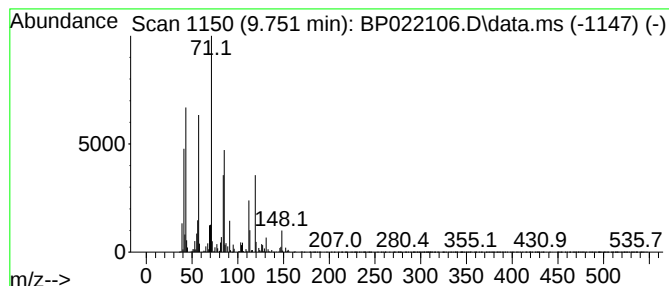
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 74

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.751	2.55 ng/ul	573819	Naphthalene-d8	10.463

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	50
2		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	43
3		Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	43
4		Sulfurous acid, 2-propyl tetra...	320	C17H36O3S	1010309-12-5	35
5		Eicosane, 10-methyl-	296	C21H44	054833-23-7	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
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Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

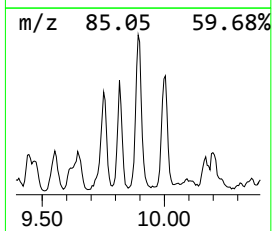
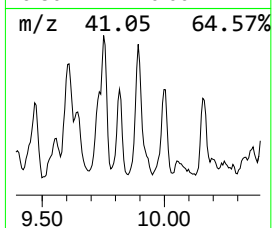
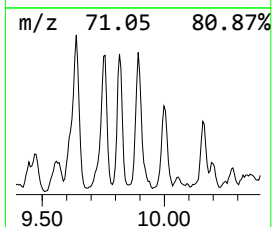
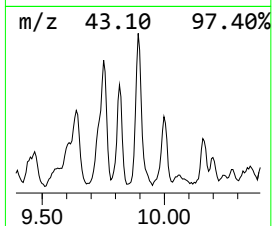
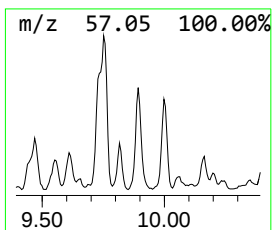
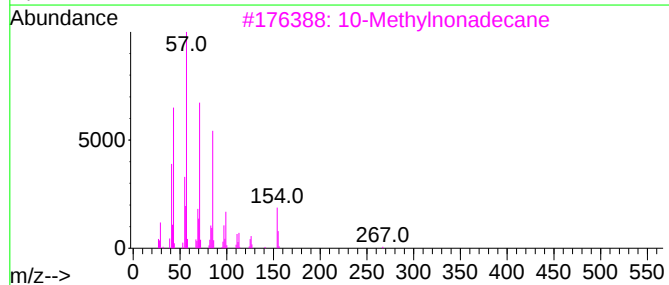
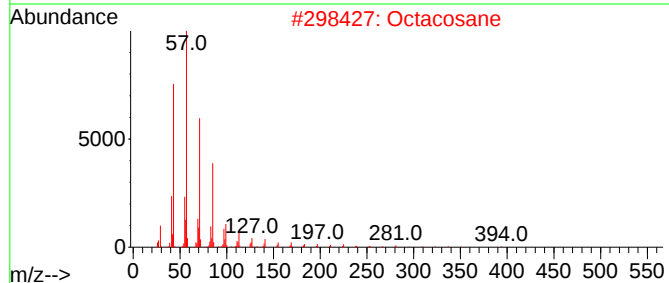
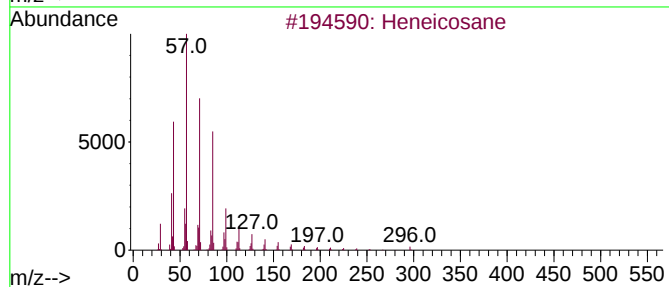
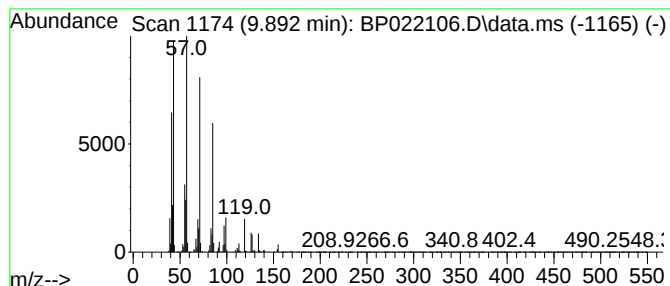
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 66

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.892	4.45 ng/ul	1001900	Naphthalene-d8	10.463	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	76
2	Octacosane	394	C28H58	000630-02-4	72
3	10-Methylnonadecane	282	C20H42	056862-62-5	72
4	Pentacosane	352	C25H52	000629-99-2	72
5	1-Undecene, 4-methyl-	168	C12H24	074630-39-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC29DL

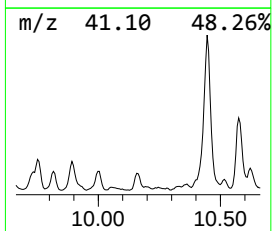
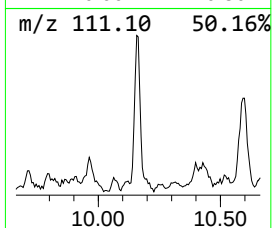
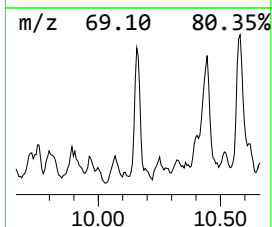
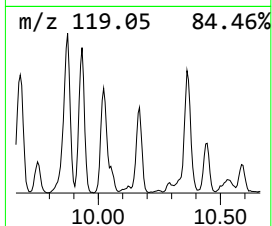
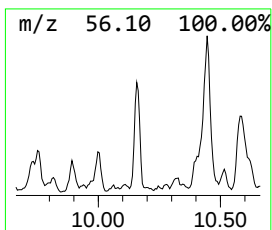
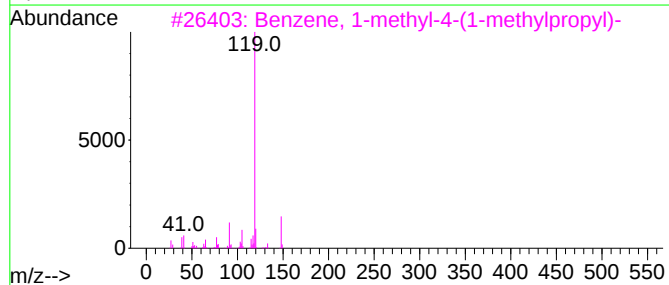
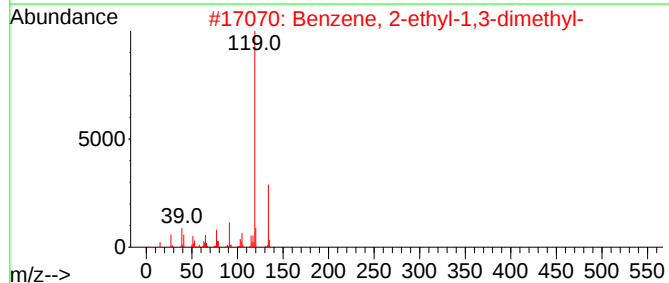
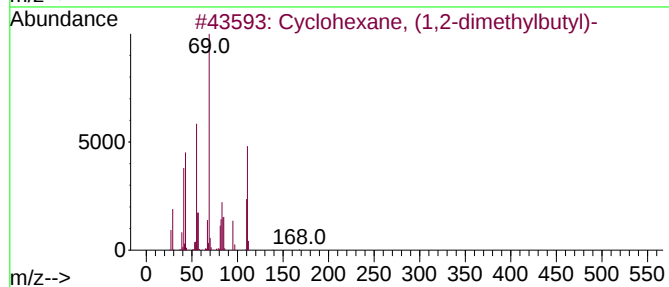
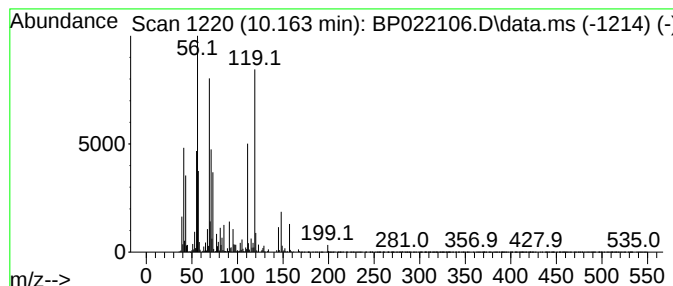
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 30 (DEL) Alkane: Cyclic10.163 Concentration Rank 75

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.163	2.31 ng/ul	519123	Naphthalene-d8	10.463

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, (1,2-dimethylbutyl)-	168	C12H24	061142-37-8	30
2		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	25
3		Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	25
4		2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	22
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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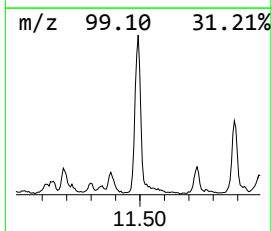
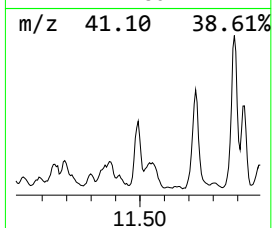
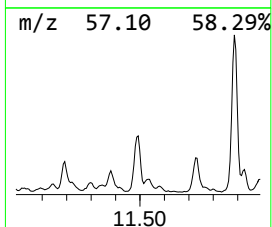
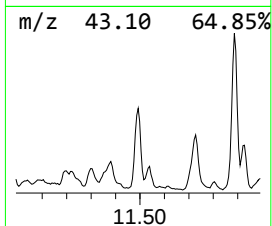
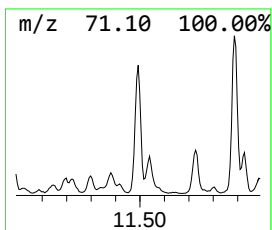
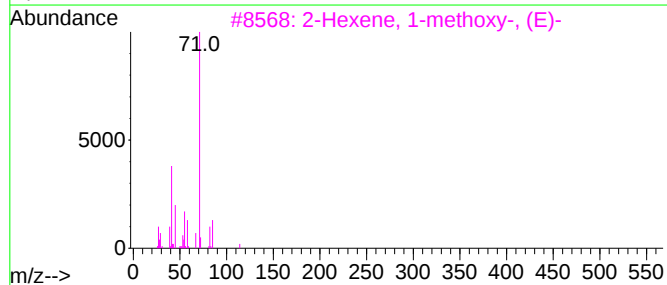
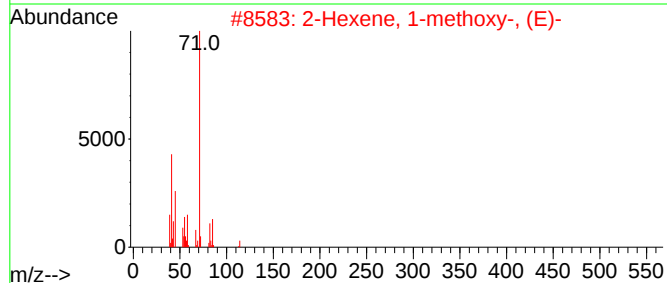
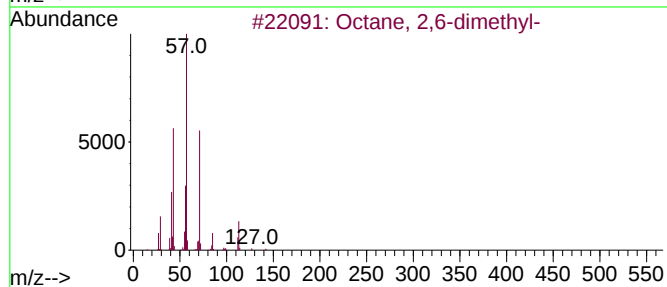
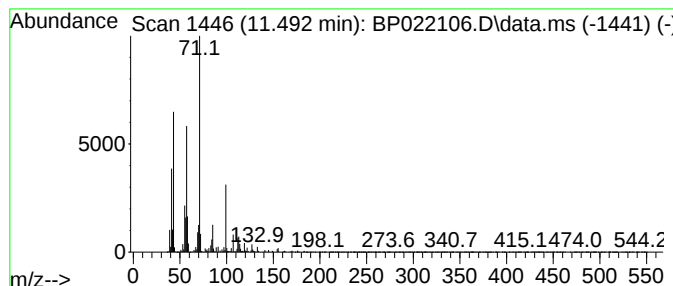
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 65

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.492	4.49 ng/ul	1010070	Naphthalene-d8	10.463

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	53	
2	2-Hexene, 1-methoxy-, (E)-	114	C7H14O	056052-83-6	52	
3	2-Hexene, 1-methoxy-, (E)-	114	C7H14O	056052-83-6	52	
4	3-Octen-2-ol	128	C8H16O	076649-14-4	52	
5	Oxalic acid, neopentyl nonyl ester	286	C16H30O4	1000309-73-3	50	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

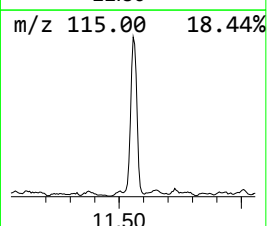
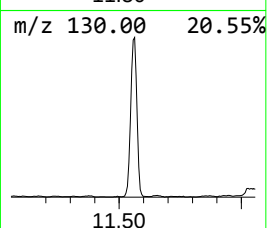
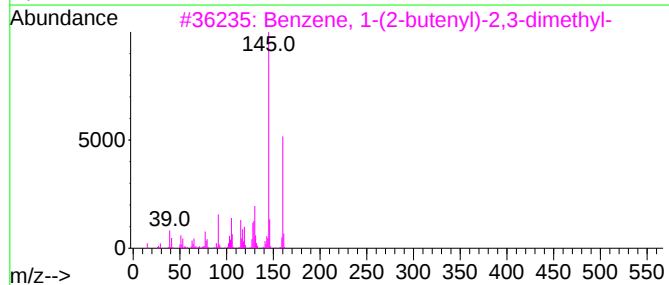
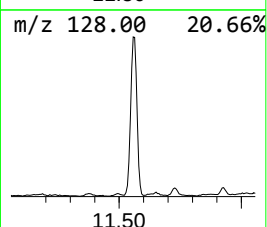
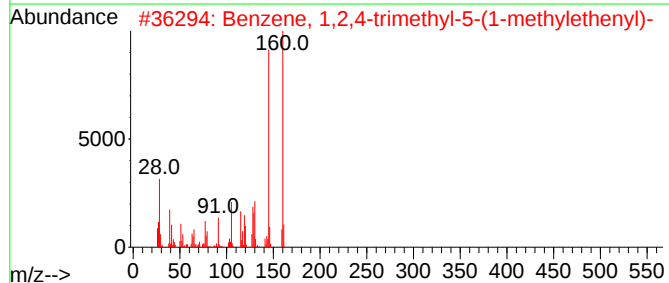
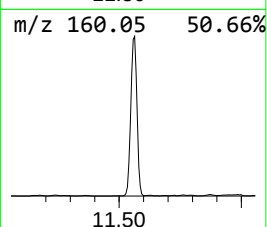
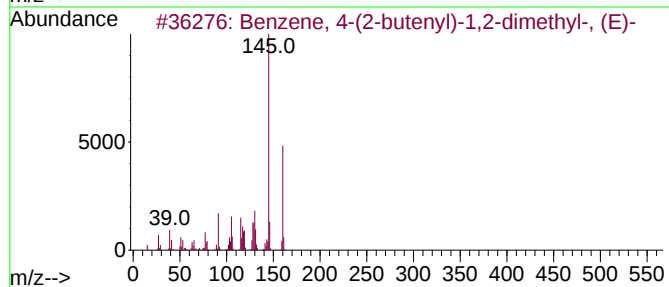
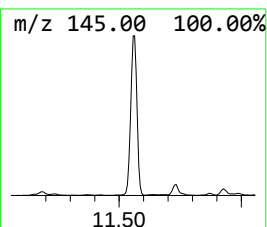
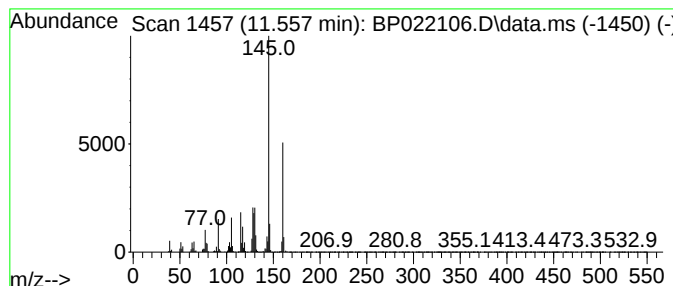
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 34 Benzene, 4-(2-butenyl)-1,2-... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.557	16.19 ng/ul	3641220	Naphthalene-d8	10.463	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	94
2	Benzene, 1,2,4-trimethyl-5-(1-me...	160	C12H16	054340-84-0	94
3	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	94
4	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	94
5	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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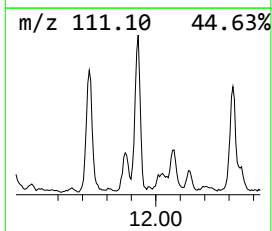
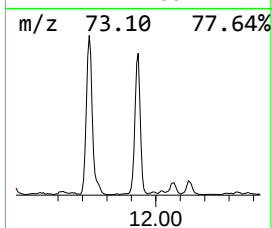
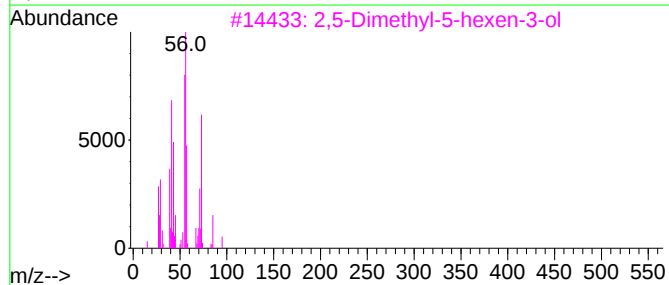
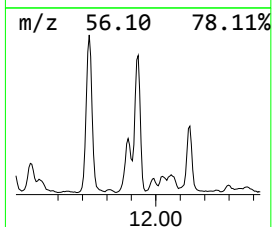
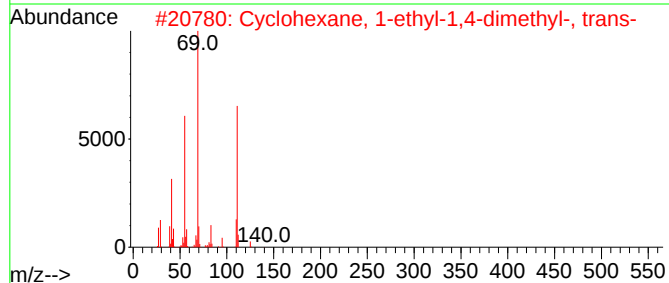
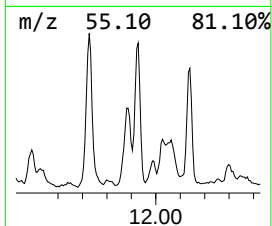
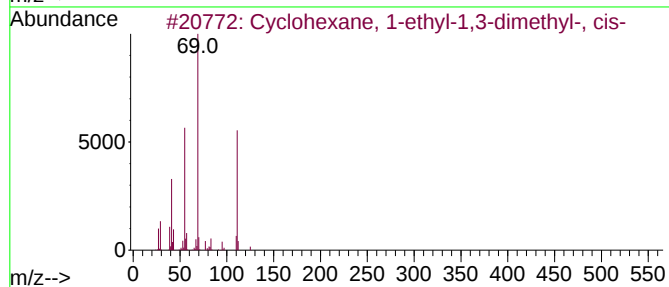
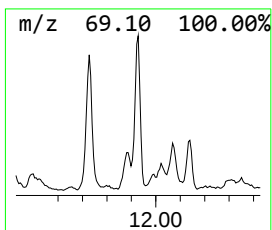
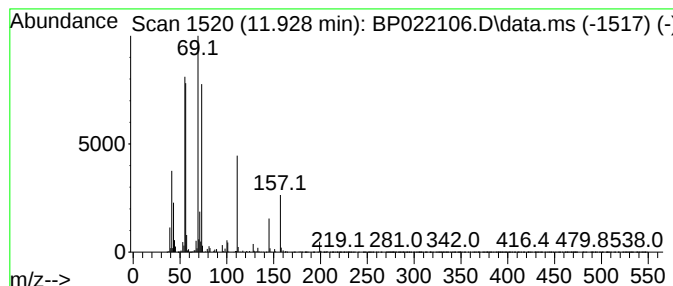
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 37 (DEL) Alkane: Cyclic11.928 Concentration Rank 64

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.928	4.77 ng/ul	1073400	Naphthalene-d8	10.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-ethyl-1,3-dimethyl...	140	C10H20	062238-31-7	35
2			Cyclohexane, 1-ethyl-1,4-dimethyl...	140	C10H20	062238-32-8	17
3			2,5-Dimethyl-5-hexen-3-ol	128	C8H16O	067760-91-2	16
4			Neopentyl glycol	104	C5H12O2	000126-30-7	12
5			Cyclopropanecarboxylic acid, 4-n...	207	C10H9NO4	1010307-52-7	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

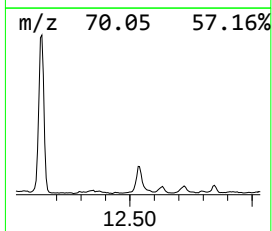
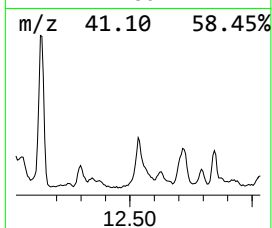
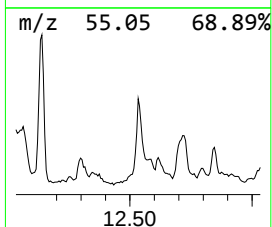
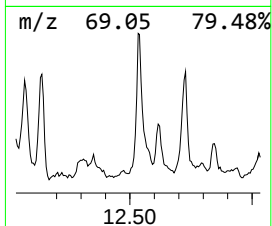
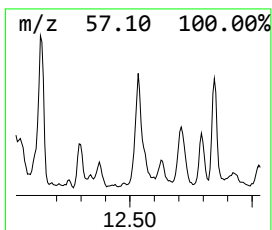
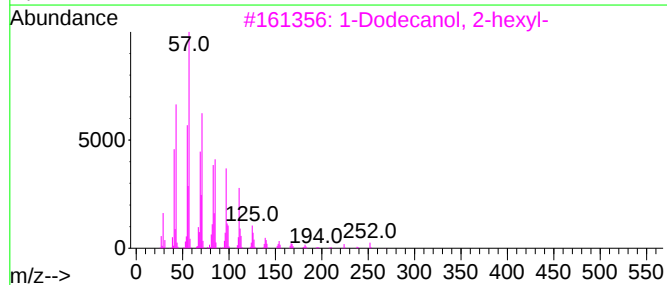
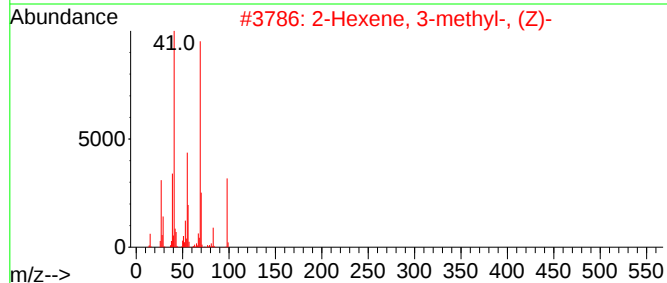
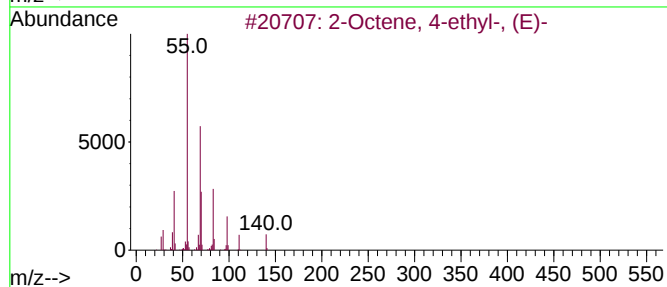
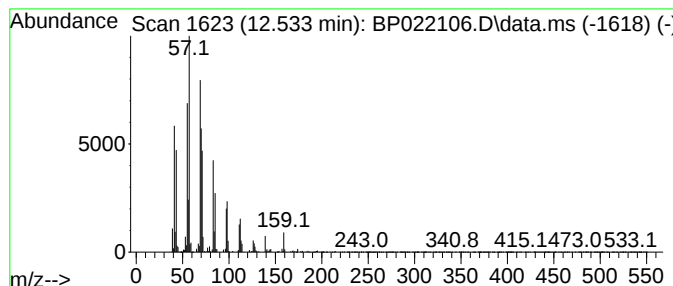
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.534	13.65 ng/ul	920632	Acenaphthene-d10		14.316	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Octene, 4-ethyl-, (E)-		140	C10H20	074630-09-4	46
2	2-Hexene, 3-methyl-, (Z)-		98	C7H14	010574-36-4	30
3	1-Dodecanol, 2-hexyl-		270	C18H38O	110225-00-8	27
4	2-Hexene, 2-methyl-		98	C7H14	002738-19-4	25
5	3-Methyl-3-hexene		98	C7H14	003404-65-7	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
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Misc :
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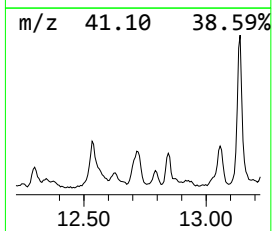
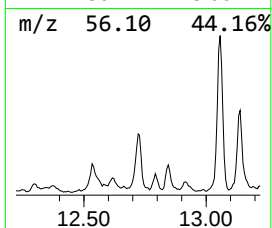
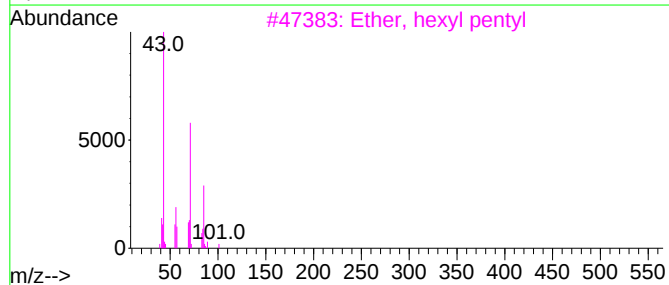
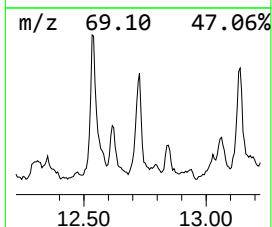
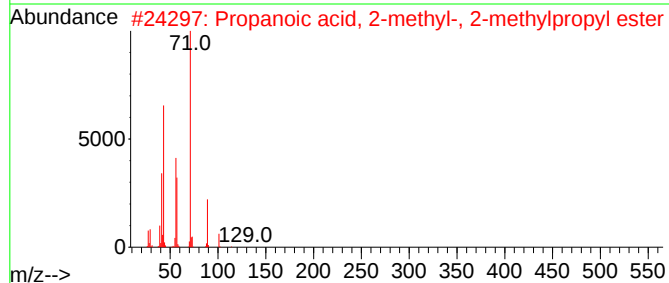
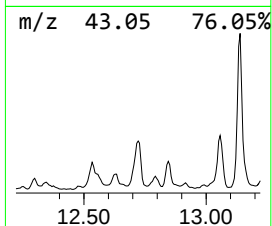
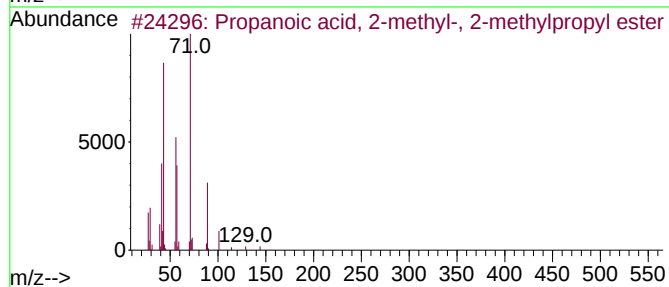
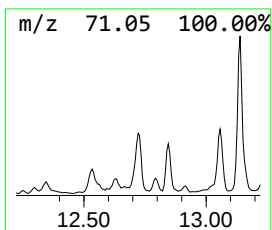
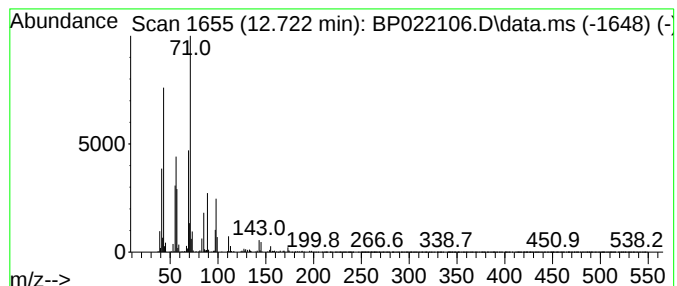
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 40 Propanoic acid, 2-methyl-, ... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.722	12.11 ng/ul	817133	Acenaphthene-d10		14.316	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-methyl-, 2-met...		144	C8H16O2	000097-85-8	53
2	Propanoic acid, 2-methyl-, 2-met...		144	C8H16O2	000097-85-8	53
3	Ether, hexyl pentyl		172	C11H24O	032357-83-8	47
4	2-Decanol, 2-methylpropionate		228	C14H28O2	1000447-30-4	47
5	2-Nonanol, 2-methylpropionate		214	C13H26O2	069121-76-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

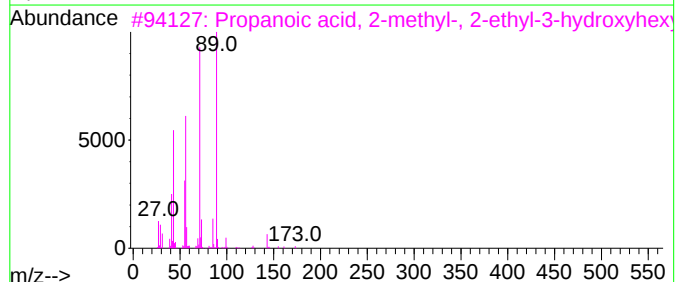
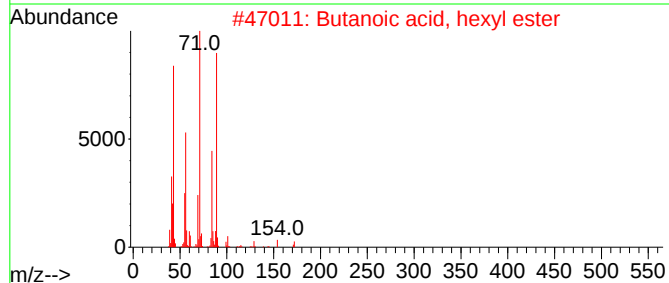
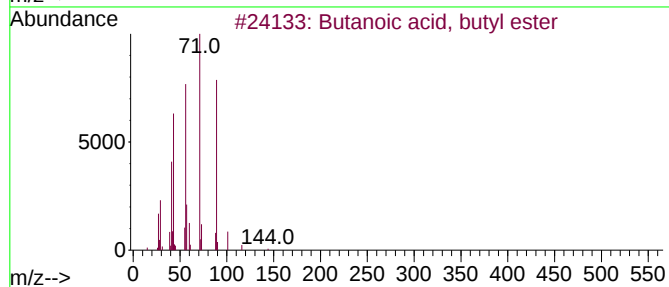
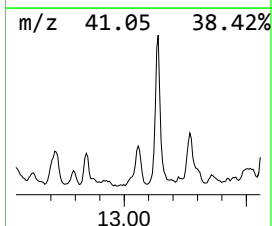
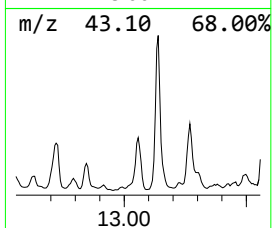
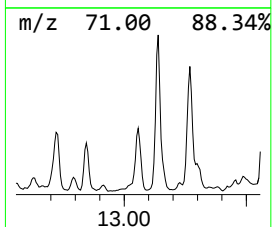
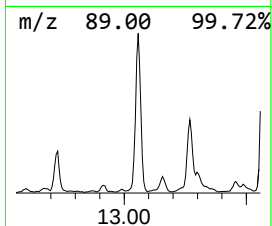
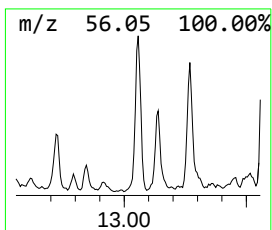
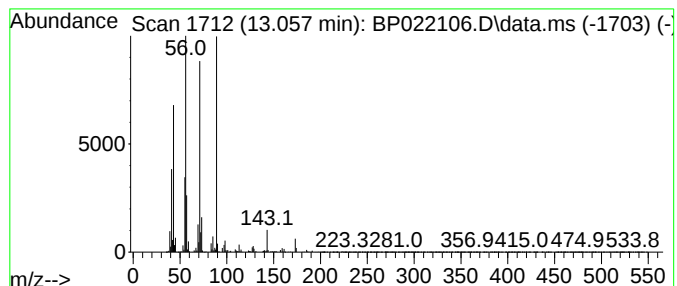
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 41 Butanoic acid, butyl ester Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.057	14.38 ng/ul	970255	Acenaphthene-d10			14.316
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Butanoic acid, butyl ester		144	C8H16O2	000109-21-7	91
2	Butanoic acid, hexyl ester		172	C10H20O2	002639-63-6	72
3	Propanoic acid, 2-methyl-, 2-eth...		216	C12H24O3	074367-31-0	64
4	Propanoic acid, 2-methyl-, 3-hyd...		216	C12H24O3	000077-68-9	64
5	Butanoic acid, butyl ester		144	C8H16O2	000109-21-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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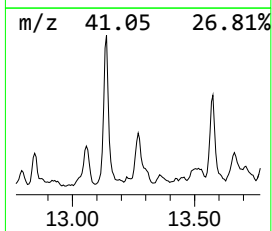
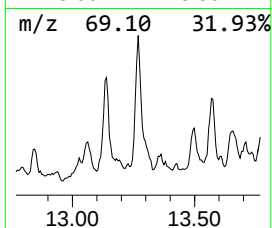
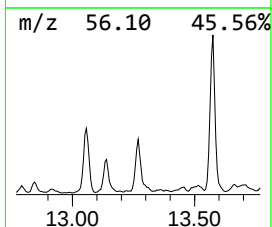
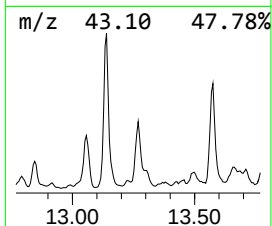
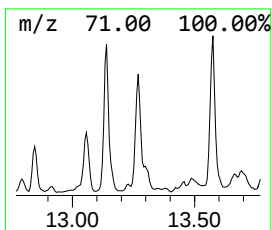
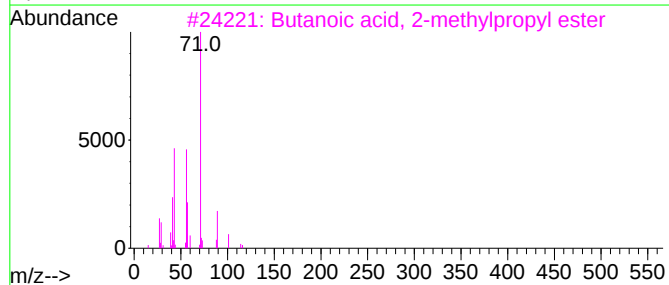
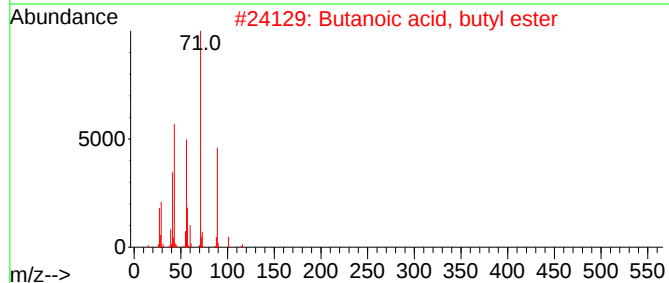
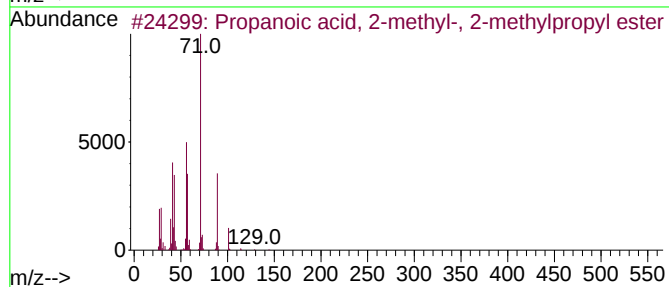
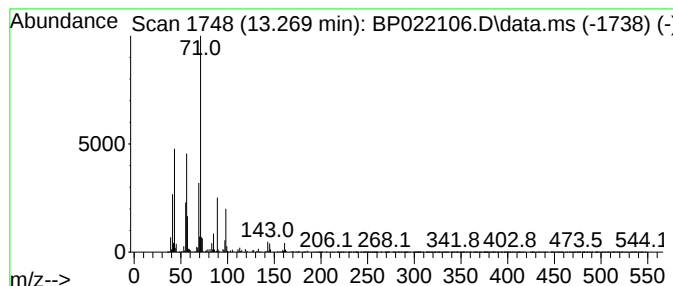
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 42 Butanoic acid, 2-methylprop... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.269	23.69 ng/ul	1598230	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	59
2			Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	59
3			Butanoic acid, 2-methylpropyl ester	144	C8H16O2	000539-90-2	59
4			Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	45
5			Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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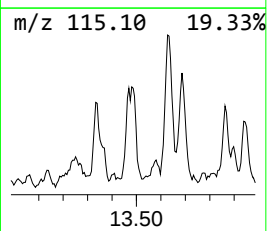
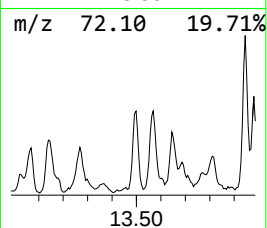
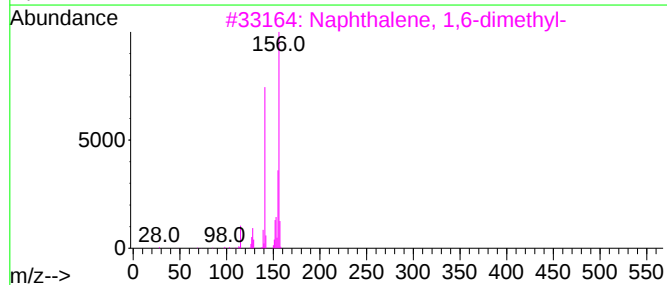
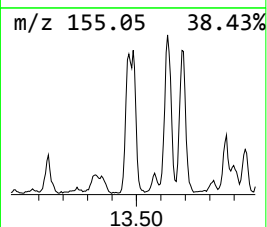
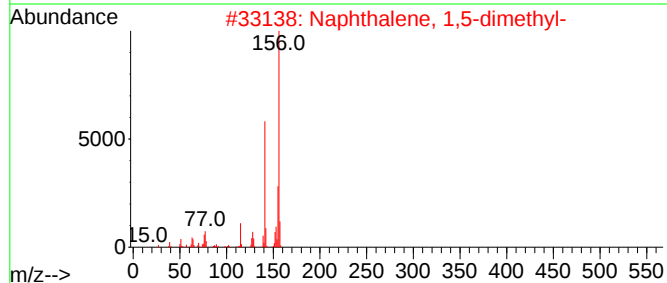
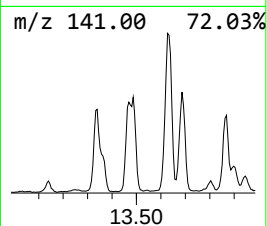
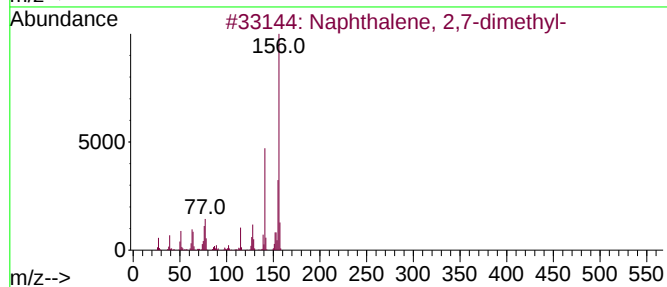
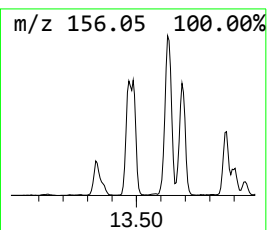
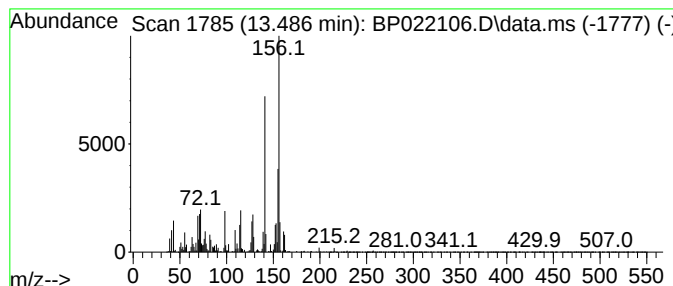
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 43 Naphthalene, 2,7-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.486	20.61 ng/ul	1390670	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
2			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	94
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	94
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	94
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

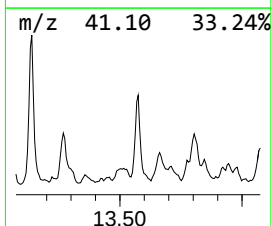
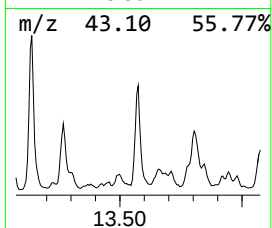
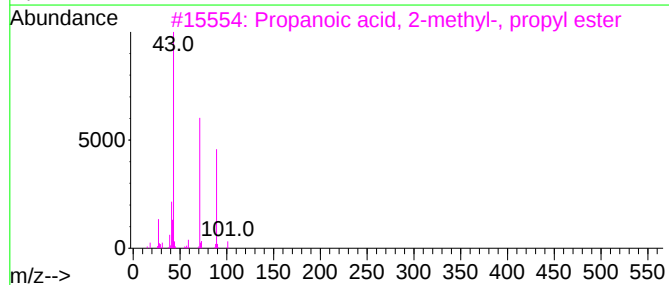
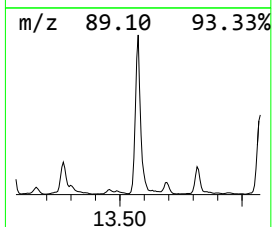
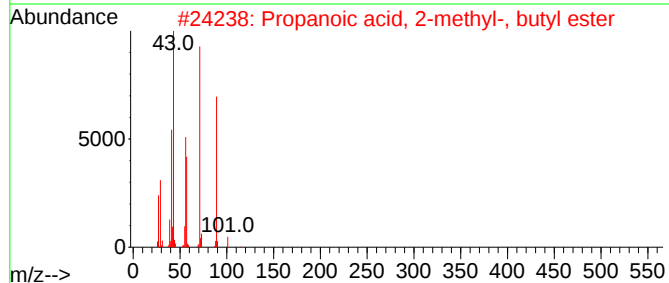
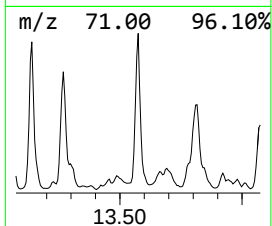
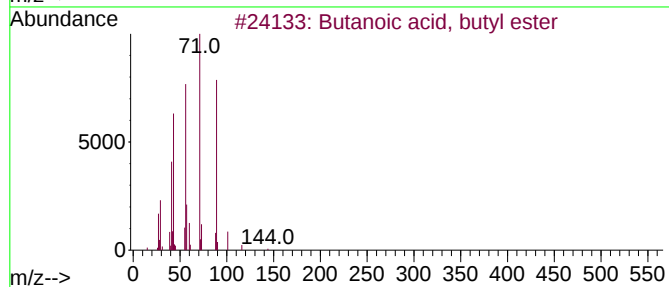
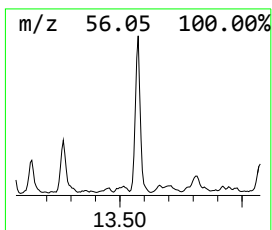
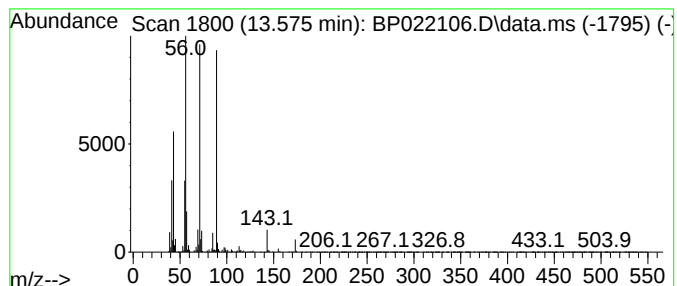
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 44 Propanoic acid, 2-methyl-, ... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.575	24.84 ng/ul	1675920	Acenaphthene-d10	14.316	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	80
2	Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	72
3	Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	59
4	Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	56
5	Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
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Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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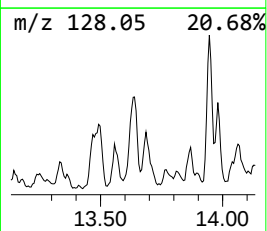
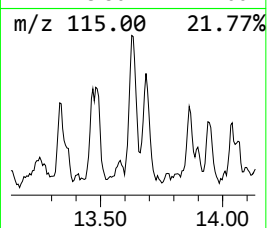
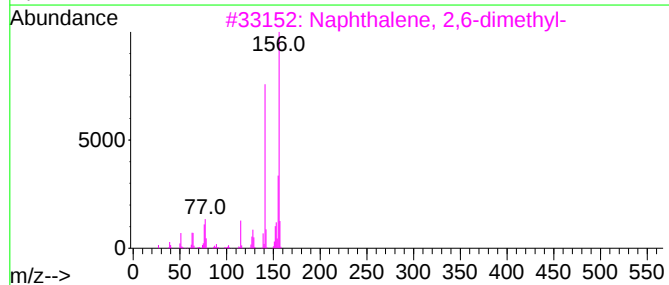
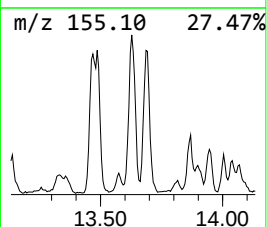
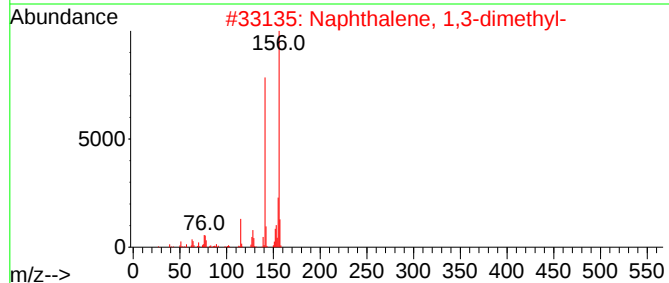
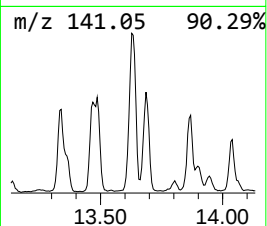
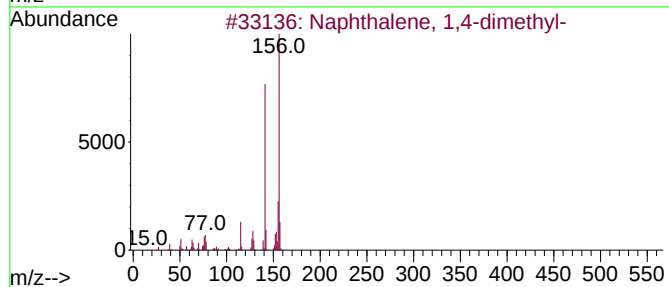
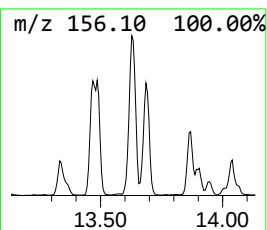
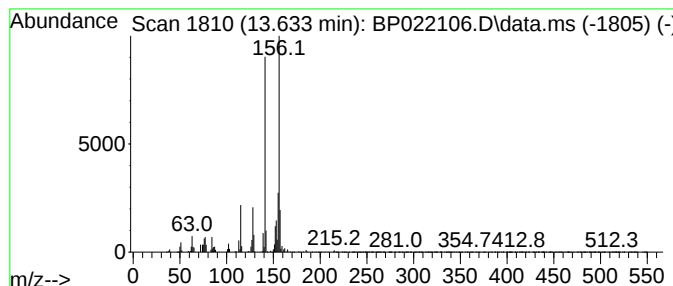
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 45 Naphthalene, 1,4-dimethyl- Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.633	9.40 ng/ul	634087	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
2			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	94
3			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	93
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	93
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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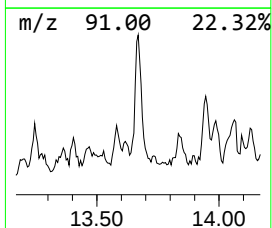
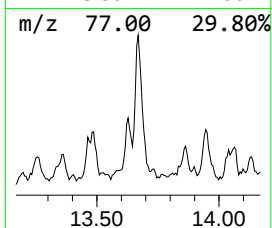
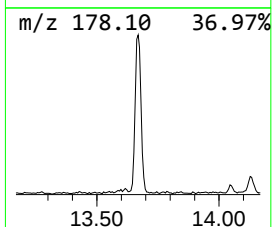
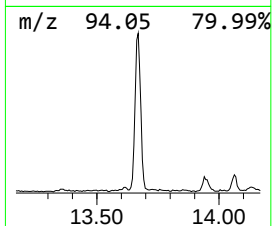
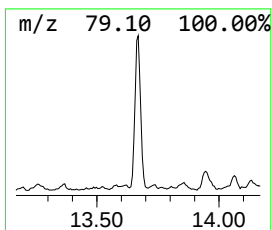
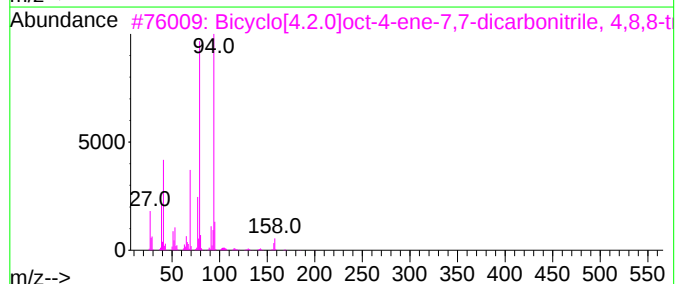
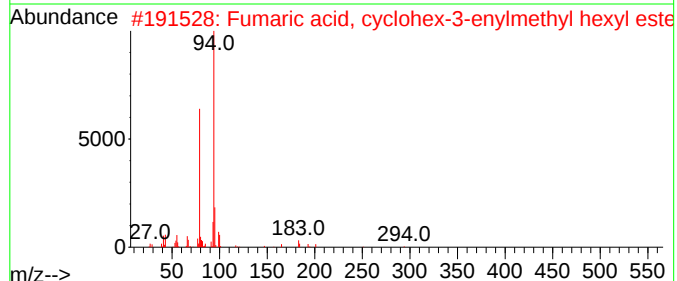
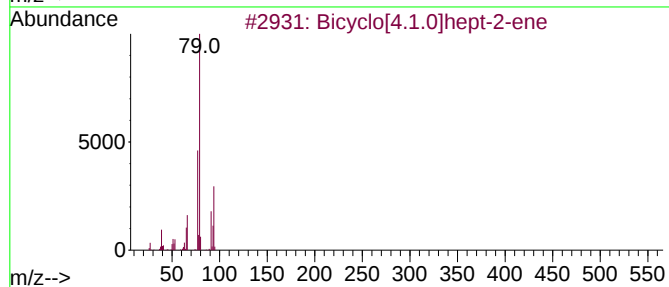
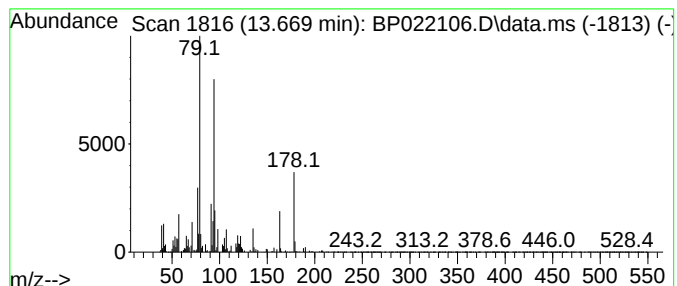
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 46 Bicyclo[4.1.0]hept-2-ene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.669	28.33 ng/ul	1911120	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[4.1.0]hept-2-ene	94	C7H10	002566-57-6	58
2			Fumaric acid, cyclohex-3-enylmet...	294	C17H26O4	1000345-14-3	53
3			Bicyclo[4.2.0]oct-4-ene-7,7-dica...	200	C13H16N2	1000163-03-5	53
4			6-Methyl-3-cyclohexen-1-carboxal...	124	C8H12O	1000132-13-3	53
5			Bicyclo[2.2.1]hept-2-ene, 2,3-di...	122	C9H14	000529-16-8	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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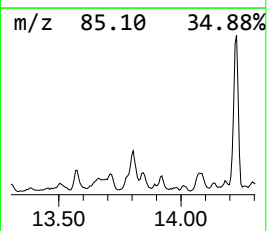
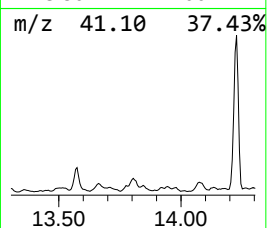
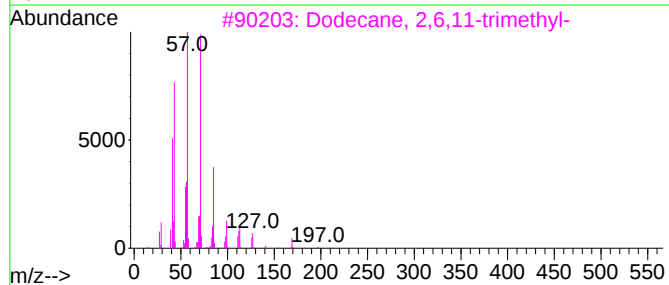
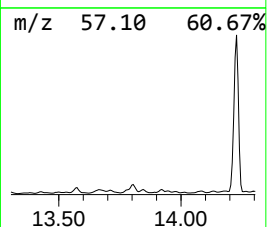
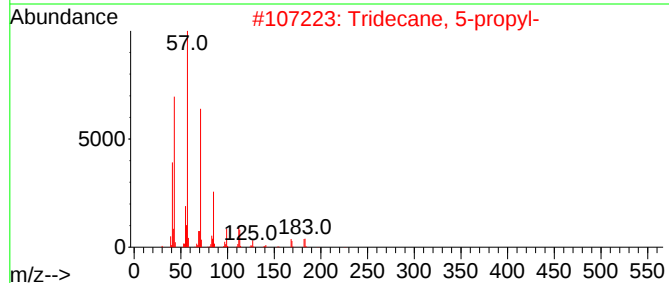
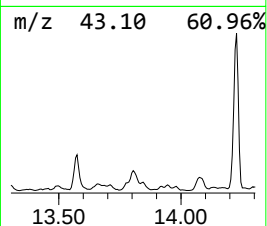
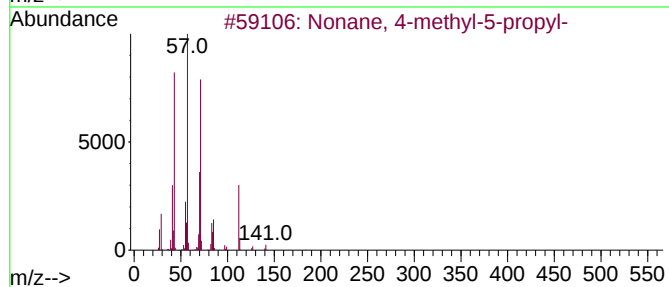
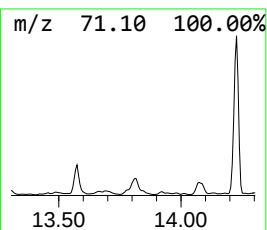
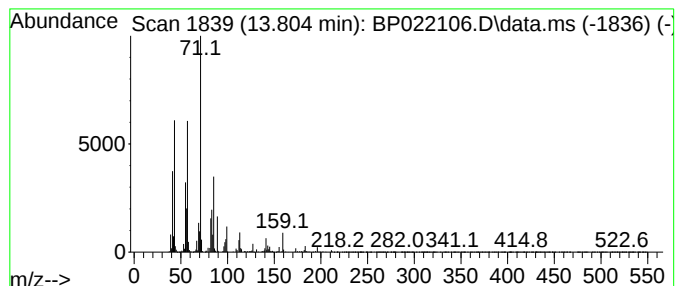
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.804	9.63 ng/ul	649577	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 4-methyl-5-propyl-	184	C13H28	062185-55-1	72
2			Tridecane, 5-propyl-	226	C16H34	055045-11-9	58
3			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	55
4			Nonadecane	268	C19H40	000629-92-5	43
5			Furazan-3-carboxamide, 4-amino-N...	212	C8H12N4O3	309735-27-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC29DL

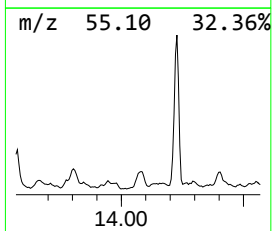
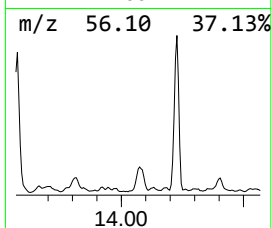
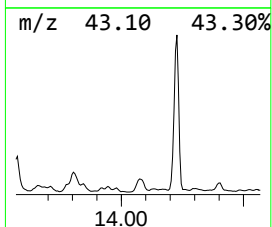
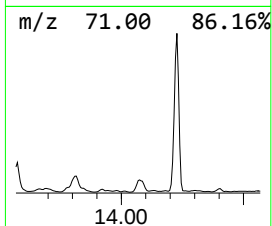
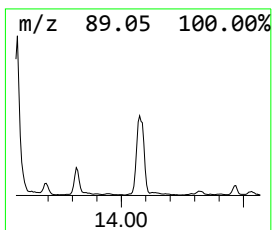
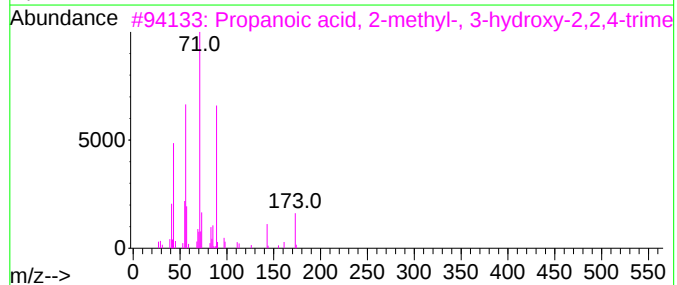
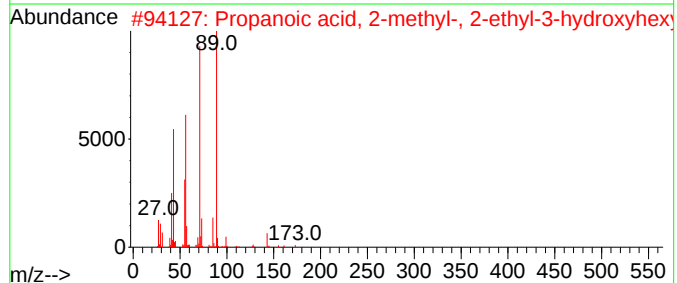
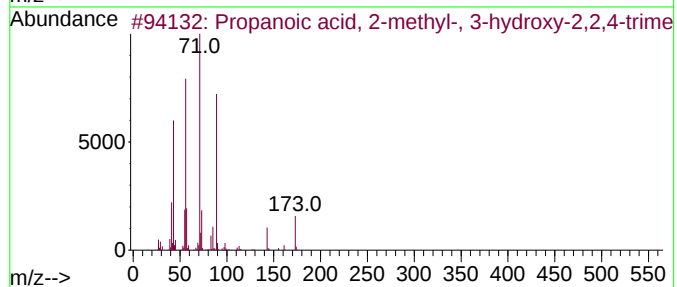
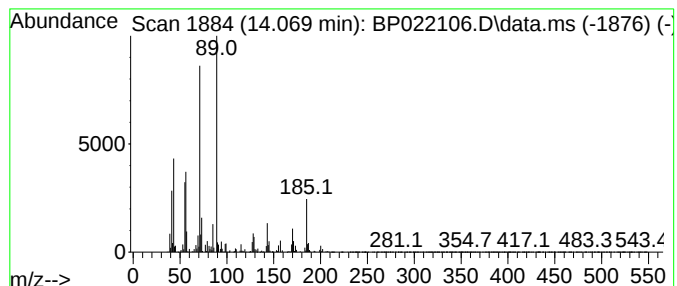
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 49 Propanoic acid, 2-methyl-, ... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.069	18.65 ng/ul	1258150	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-methyl-, 3-hyd...	216	C12H24O3	000077-68-9	78
2			Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	78
3			Propanoic acid, 2-methyl-, 3-hyd...	216	C12H24O3	000077-68-9	64
4			Butanoic acid, decyl ester	228	C14H28O2	005454-09-1	56
5			Butyric acid, tetradecyl ester	284	C18H36O2	006221-98-3	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC29DL

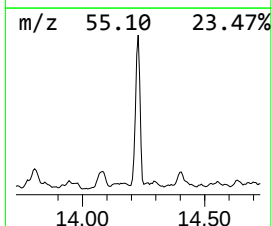
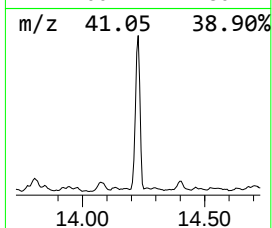
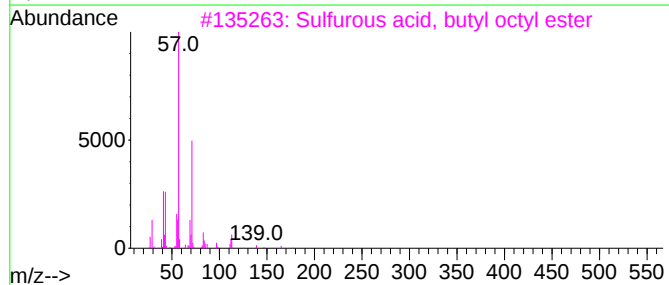
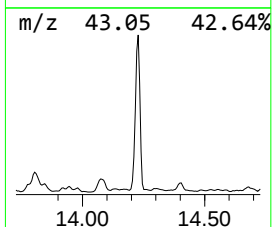
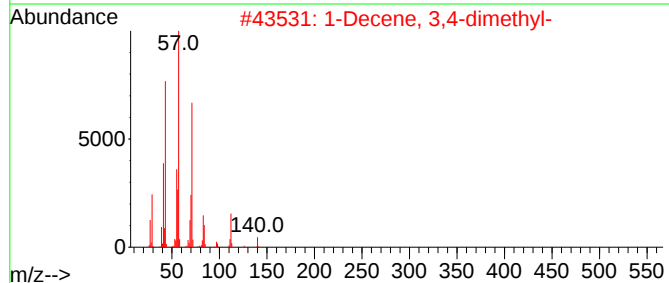
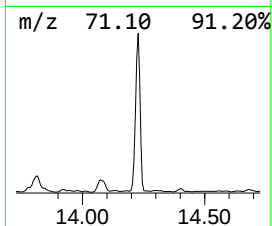
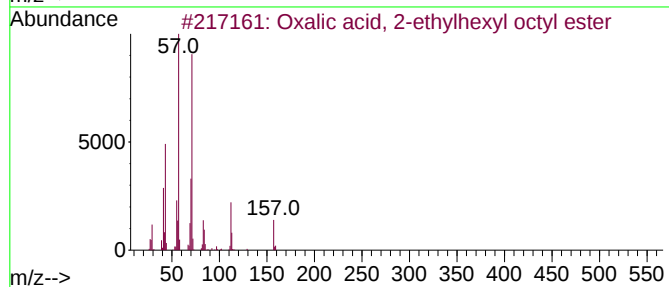
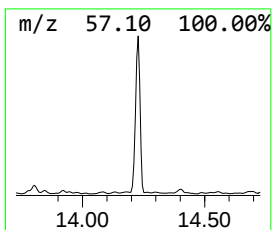
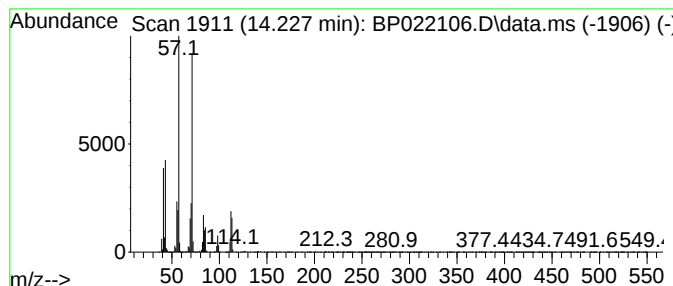
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 50 Oxalic acid, 2-ethylhexyl o... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.227	119.00 ng/ul	8027580	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, 2-ethylhexyl octyl ...	314	C18H34O4	1000309-39-1	80
2			1-Decene, 3,4-dimethyl-	168	C12H24	050871-03-9	78
3			Sulfurous acid, butyl octyl ester	250	C12H26O3S	1000309-17-5	72
4			Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	72
5			Decyl pentyl ether	228	C15H32O	1000406-41-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

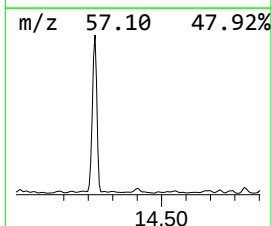
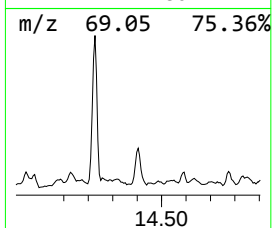
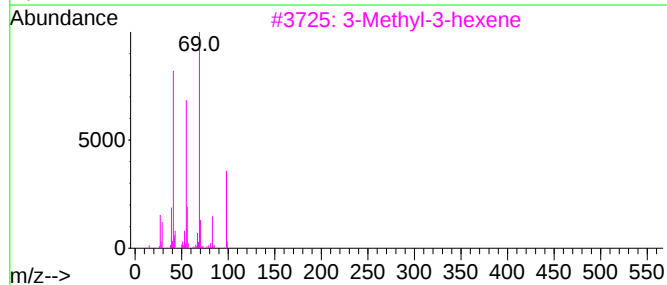
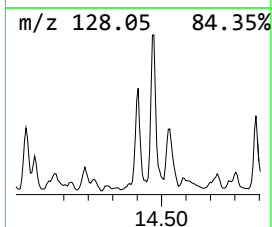
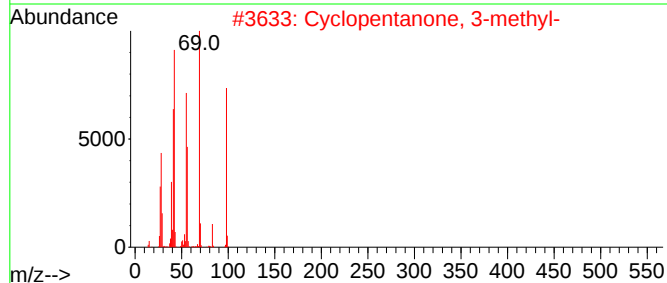
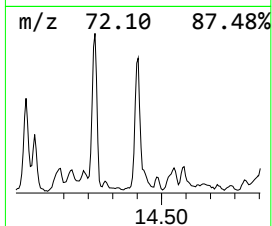
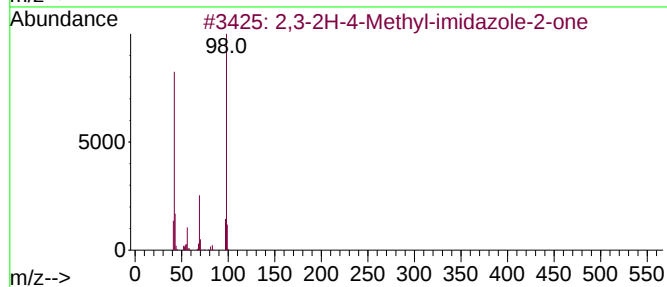
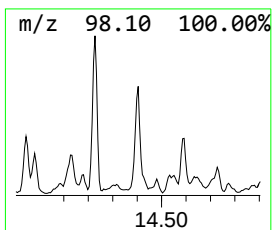
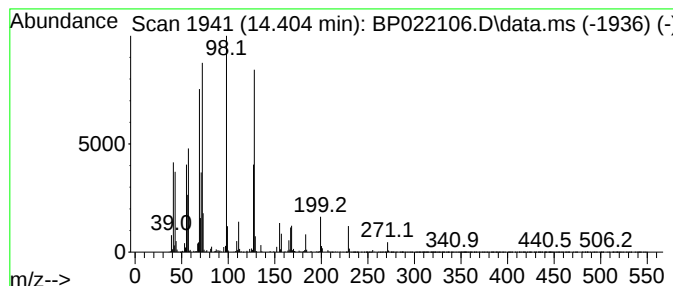
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 51 unknown-02 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.404	15.65 ng/ul	1055650	Acenaphthene-d10	14.316	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3-2H-4-Methyl-imidazole-2-one	98	C4H6N2O	039799-77-4	27
2	Cyclopentanone, 3-methyl-	98	C6H10O	001757-42-2	27
3	3-Methyl-3-hexene	98	C7H14	003404-65-7	27
4	(R)-(+)-3-Methylcyclopentanone	98	C6H10O	006672-30-6	27
5	Cyclopropanecarboxamide, N-propy...	183	C11H21NO	1000437-75-3	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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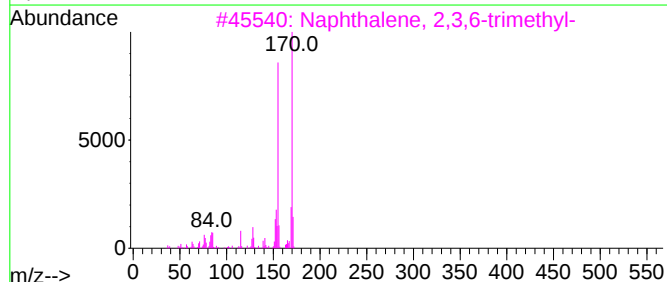
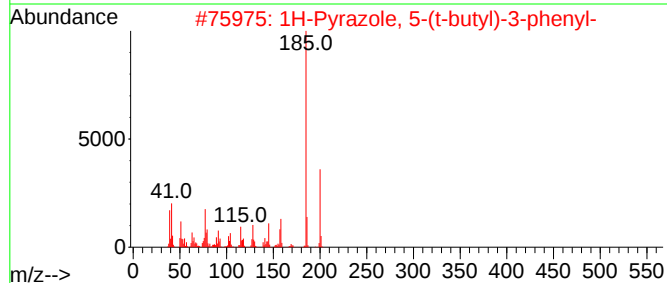
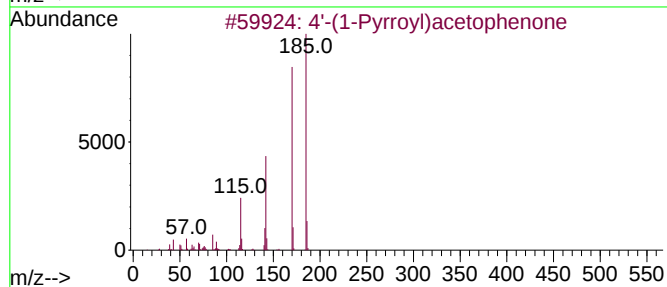
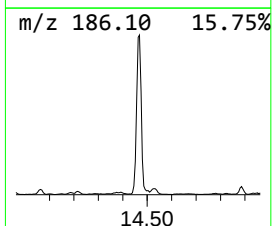
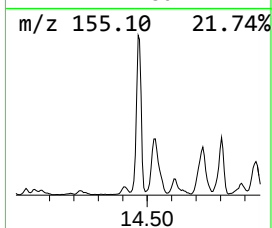
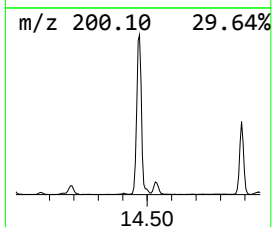
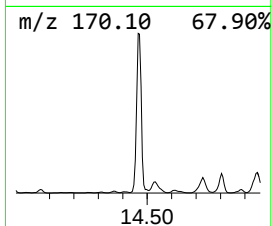
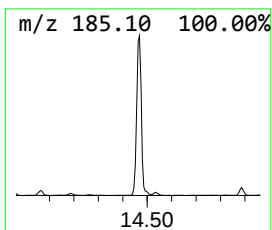
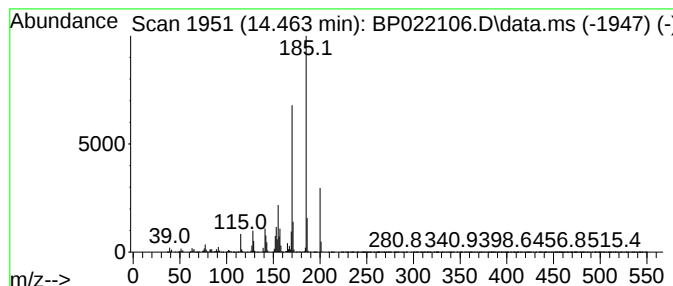
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 52 4'-(1-Pyrrolyl)acetophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.463	70.48 ng/ul	4754240	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4'-(1-Pyrrolyl)acetophenone	185	C12H11NO	022106-37-2	58
2			1H-Pyrazole, 5-(t-butyl)-3-phenyl-	200	C13H16N2	1000261-34-8	55
3			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	51
4			Cyclopropane, 1-methoxy-2,2-dime...	200	C14H16O	1000271-83-0	50
5			6-Fluoro-2,4-pyrimidinediamine, TMS	200	C7H13FN4Si	1000472-66-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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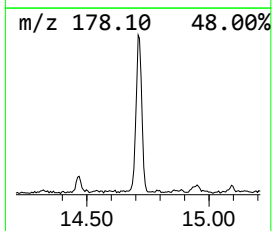
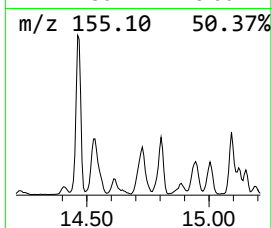
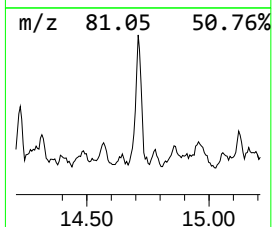
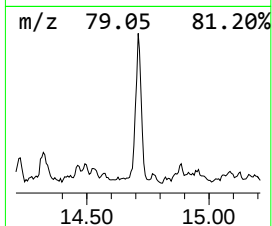
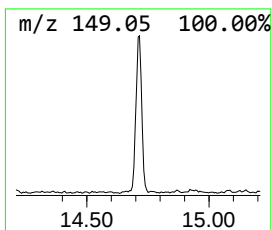
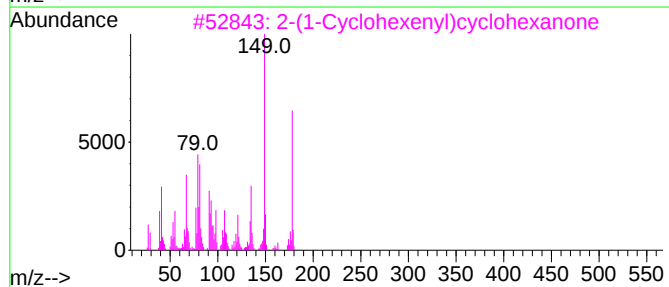
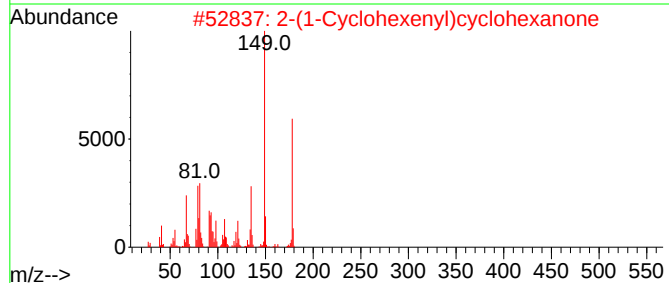
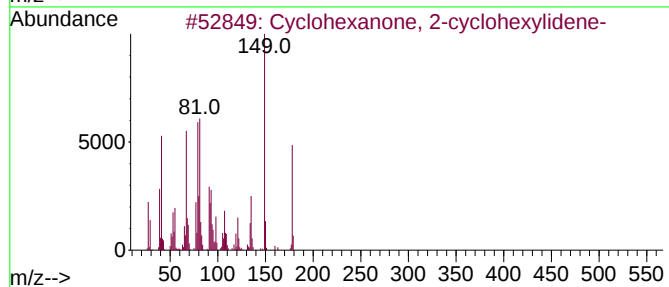
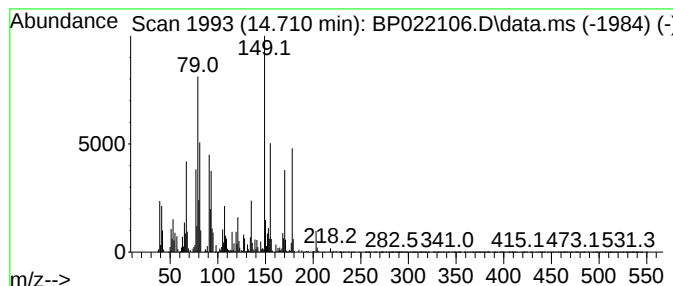
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 53 Cyclohexanone, 2-cyclohexyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.710	26.30 ng/ul	1774300	Acenaphthene-d10	14.316

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexanone, 2-cyclohexylidene-	178	C12H18O	001011-12-7	83
2		2-(1-Cyclohexenyl)cyclohexanone	178	C12H18O	001502-22-3	60
3		2-(1-Cyclohexenyl)cyclohexanone	178	C12H18O	001502-22-3	49
4		5-Pentylcyclohexa-1,3-diene	150	C11H18	056318-84-4	38
5		Cyclohexanone, 2-cyclohexylidene-	178	C12H18O	001011-12-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
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Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

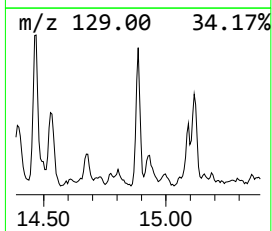
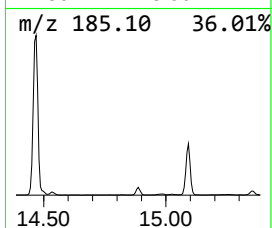
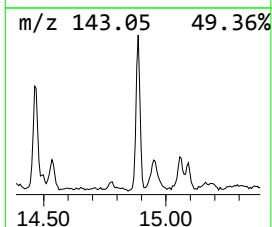
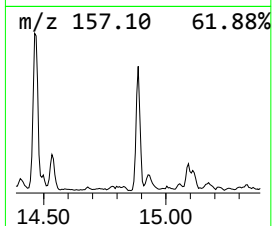
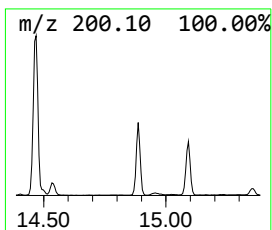
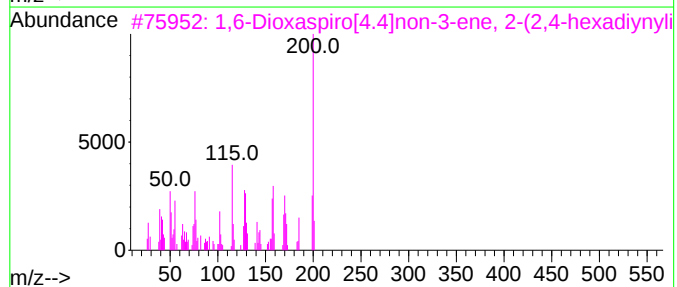
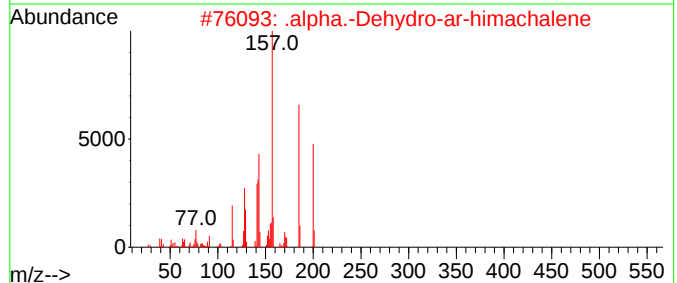
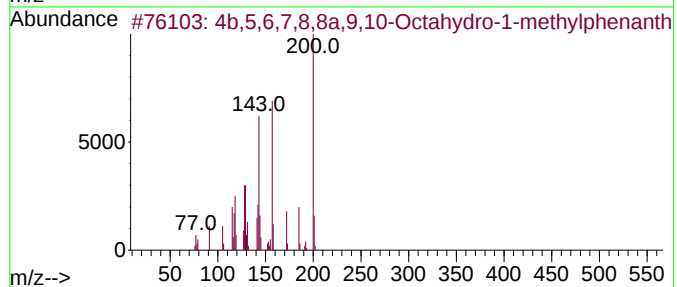
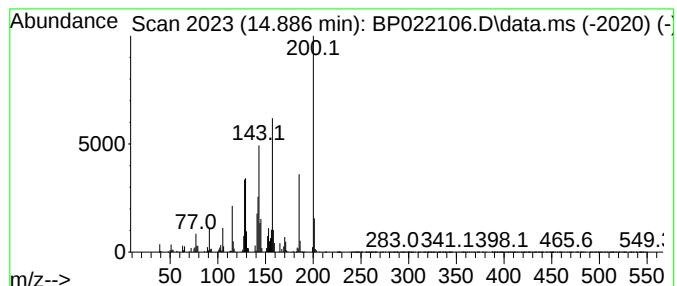
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 54 4b,5,6,7,8,8a,9,10-Octahydr... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.886	13.07 ng/ul	881643	Acenaphthene-d10	14.316	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4b,5,6,7,8,8a,9,10-Octahydro-1-m...	200	C15H20	1000080-19-3	91
2	.alpha.-Dehydro-ar-himachalene	200	C15H20	078204-62-3	64
3	1,6-Dioxaspiro[4.4]non-3-ene, 2-...	200	C13H12O2	016863-61-9	52
4	Anthracene, 1,2,3,4,5,6,7,8-octa...	200	C15H20	1000159-30-4	52
5	2-(4-Hydroxy-3-methoxybenzyliden...	200	C11H8N2O2	003696-12-6	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC29DL

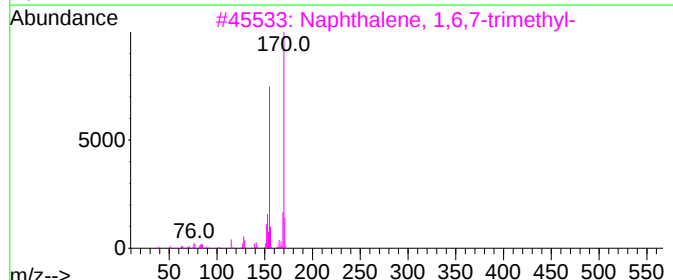
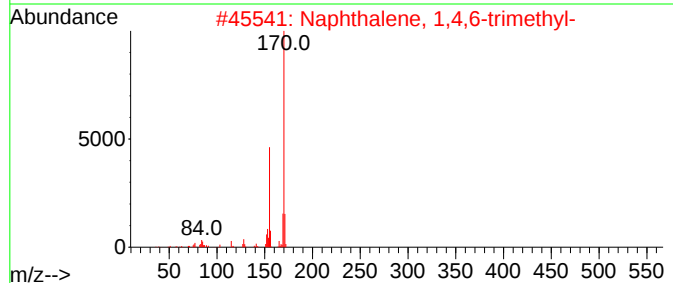
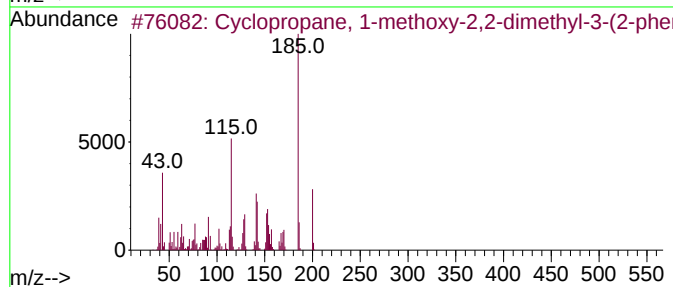
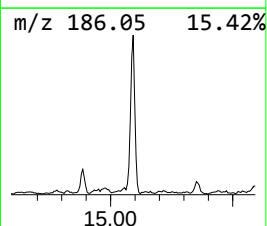
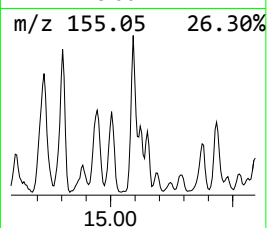
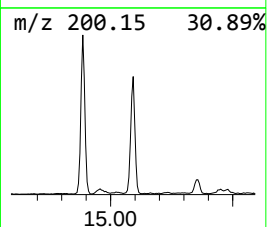
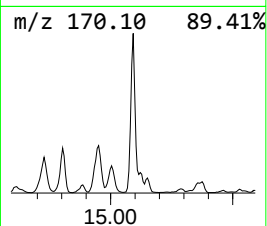
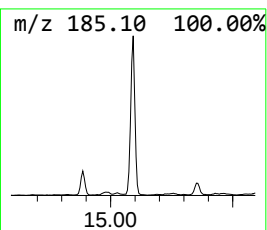
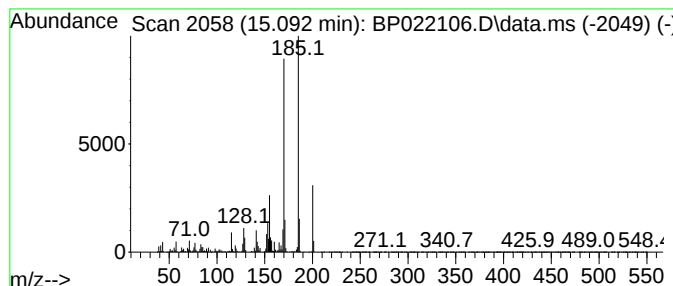
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 55 Cyclopropane, 1-methoxy-2,2... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.092	33.64 ng/ul	2269180	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1-methoxy-2,2-dime...	200	C14H16O	1000271-83-0	55
2			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	50
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	46
4			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	46
5			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC29DL

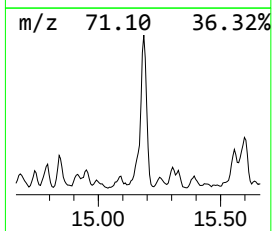
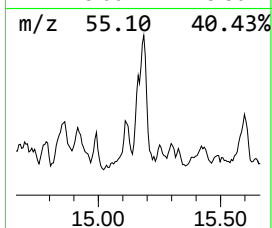
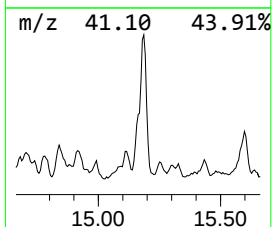
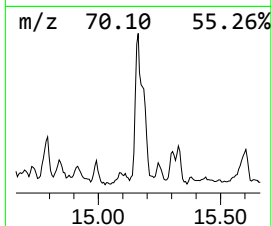
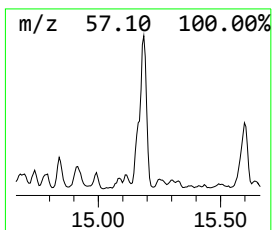
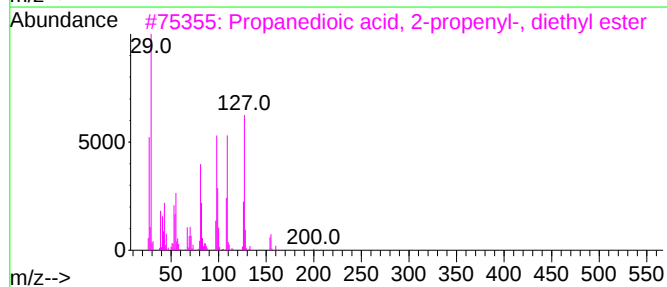
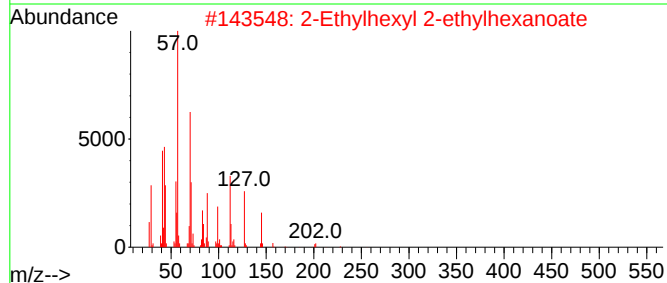
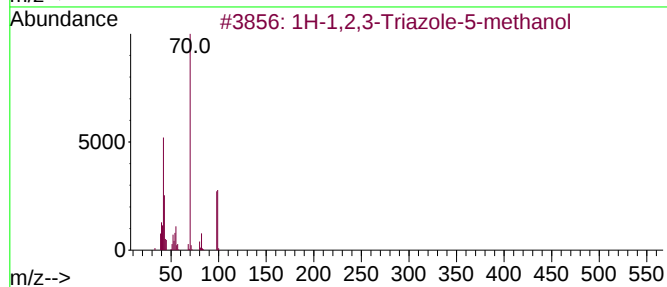
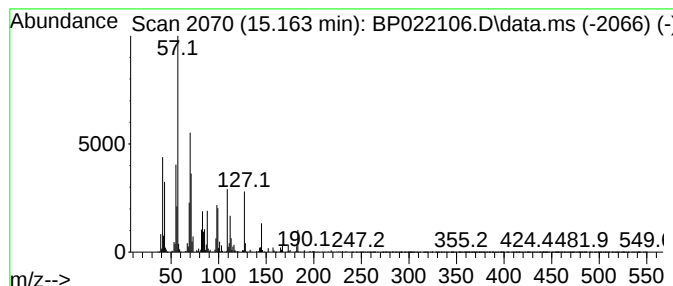
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 56 unknown-03 Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.163	11.34 ng/ul	765031	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-1,2,3-Triazole-5-methanol	99	C3H5N3O	1010319-26-5	35
2			2-Ethylhexyl 2-ethylhexanoate	256	C16H32O2	007425-14-1	27
3			Propanedioic acid, 2-propenyl-, ...	200	C10H16O4	002049-80-1	25
4			2-Ethylhexyl 2-ethylhexanoate	256	C16H32O2	007425-14-1	22
5			2-Methylbutyl octanoate	214	C13H26O2	067121-39-5	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

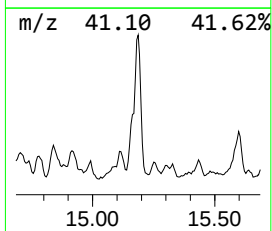
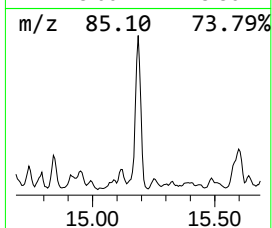
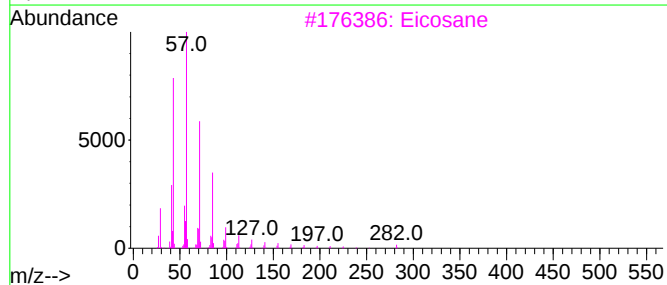
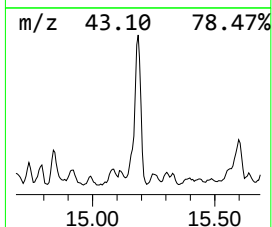
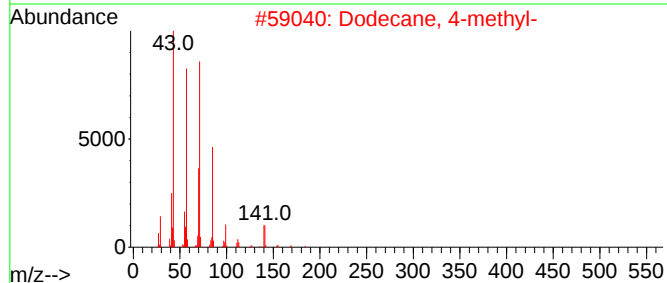
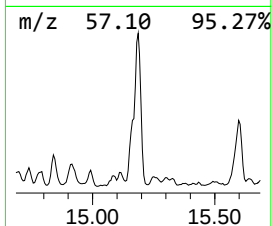
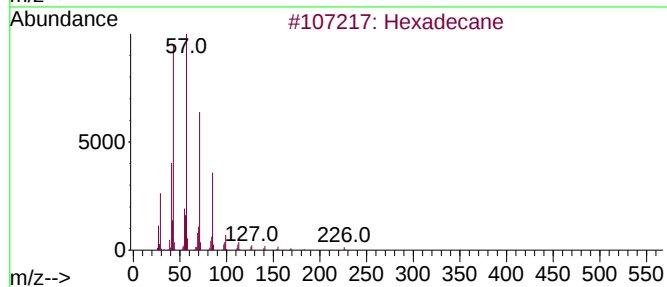
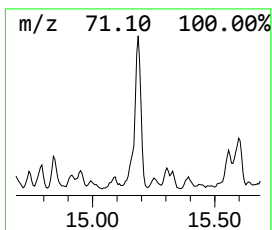
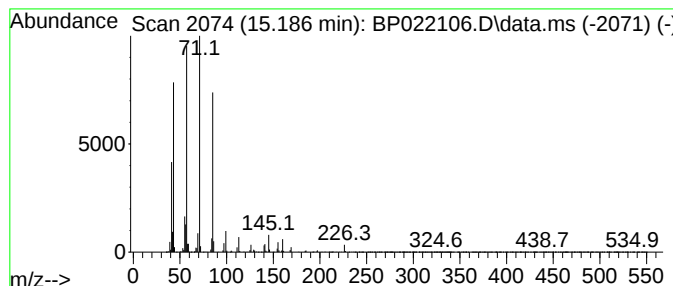
Instrument :
BNA_P
ClientSampleId :
DDC29DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 57 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.186	28.20 ng/ul	1902440	Acenaphthene-d10	14.316	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	80
2	Dodecane, 4-methyl-	184	C13H28	006117-97-1	58
3	Eicosane	282	C20H42	000112-95-8	43
4	Hexadecane	226	C16H34	000544-76-3	42
5	Hexadecane	226	C16H34	000544-76-3	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

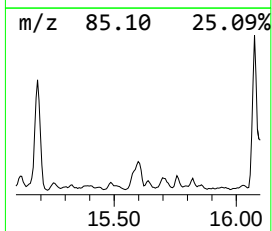
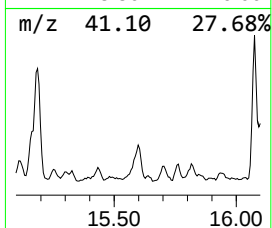
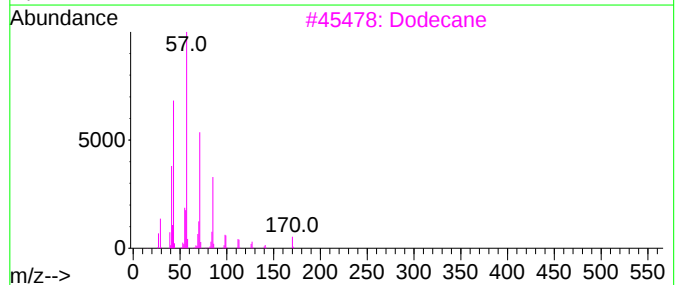
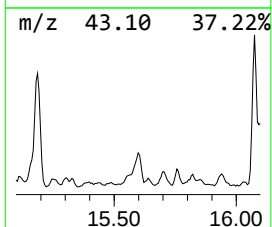
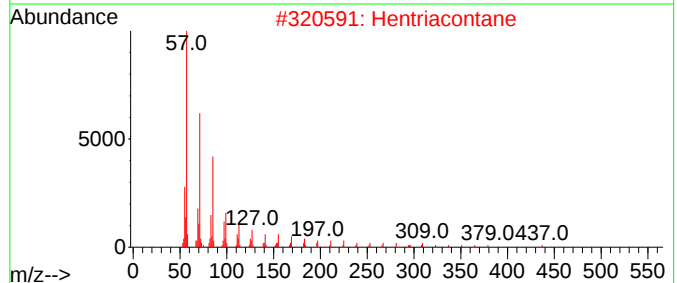
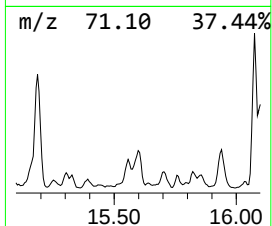
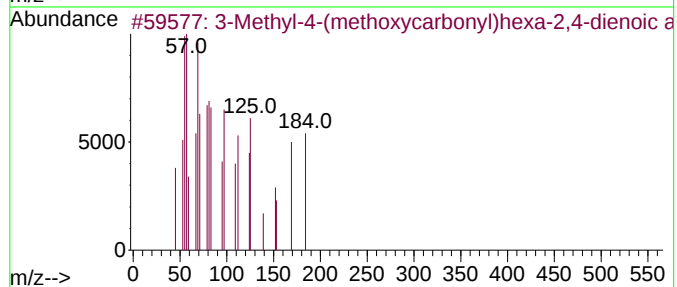
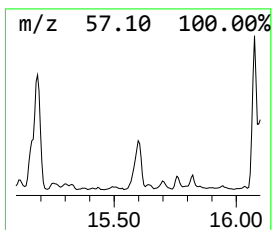
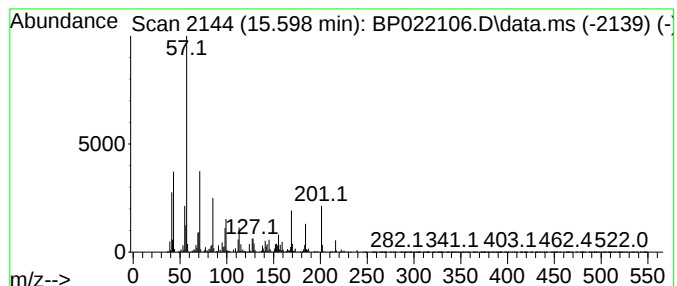
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 59 3-Methyl-4-(methoxycarbonyl... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.598	14.23 ng/ul	959929	Acenaphthene-d10	14.316	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methyl-4-(methoxycarbonyl)hexa...	184	C9H12O4	1010104-10-8	74
2	Hentriacontane	437	C31H64	000630-04-6	38
3	Dodecane	170	C12H26	000112-40-3	38
4	Hexadecane	226	C16H34	000544-76-3	35
5	Undecane, 5-ethyl-	184	C13H28	017453-94-0	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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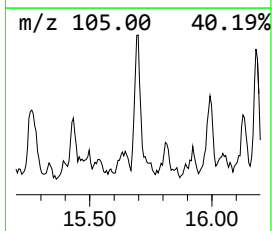
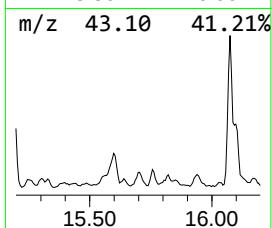
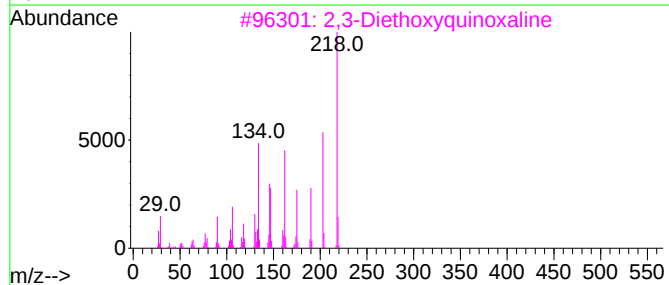
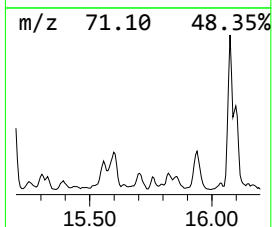
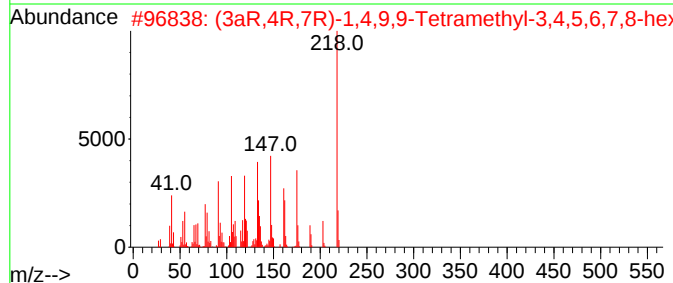
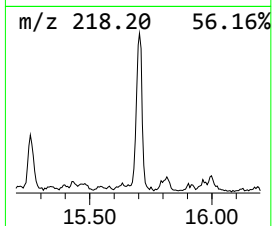
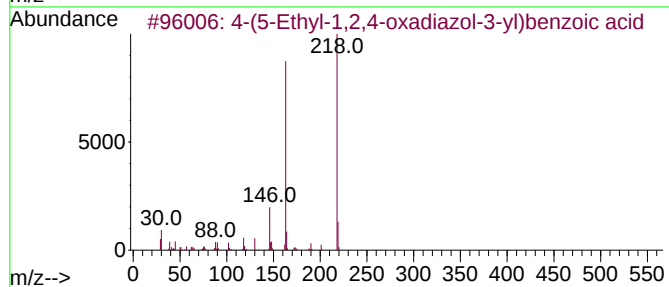
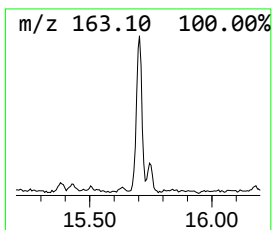
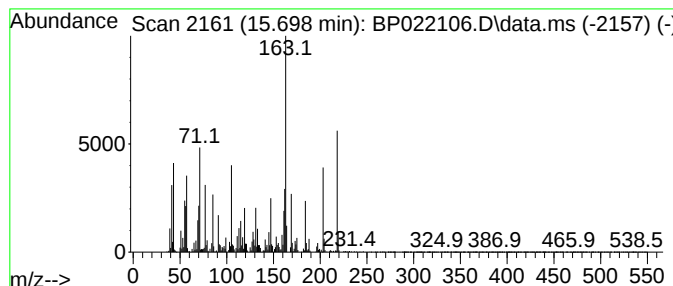
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 60 unknown-04 Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.698	12.12 ng/ul	817679	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(5-Ethyl-1,2,4-oxadiazol-3-yl)...	218	C11H10N2O3	769132-76-5	27		
2	(3aR,4R,7R)-1,4,9,9-Tetramethyl-...	218	C15H22O	003466-15-7	25		
3	2,3-Diethoxyquinoxaline	218	C12H14N2O2	057050-66-5	22		
4	1H-isoindole-1,3(2H)-dione, 5-hy...	163	C8H5N3O2	1000404-43-3	14		
5	5-Methyl-2-phenyl-2-propyle-4-he...	232	C16H24O	1000432-45-6	14		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
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Instrument :
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ClientSampleId :
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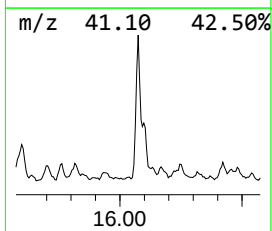
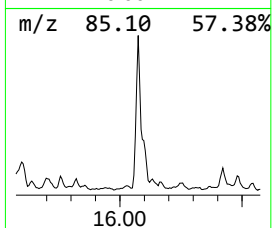
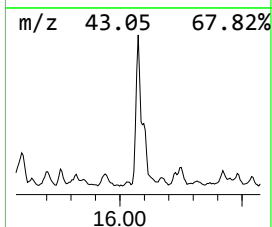
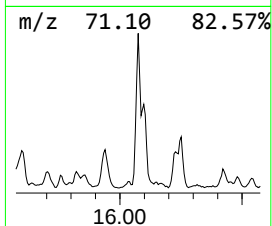
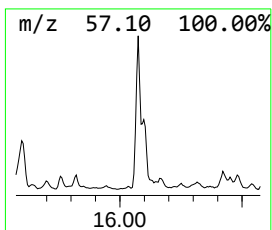
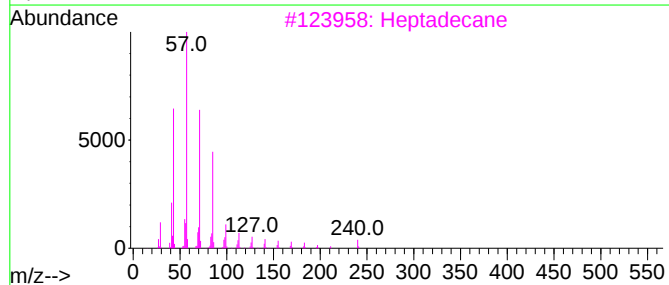
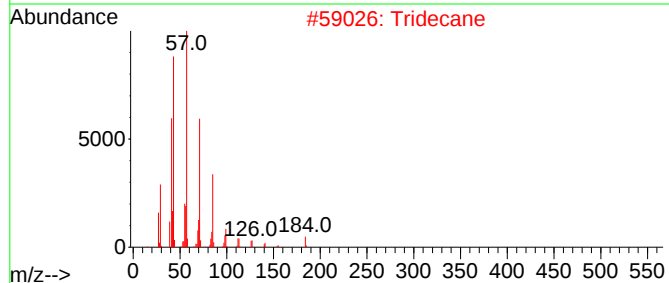
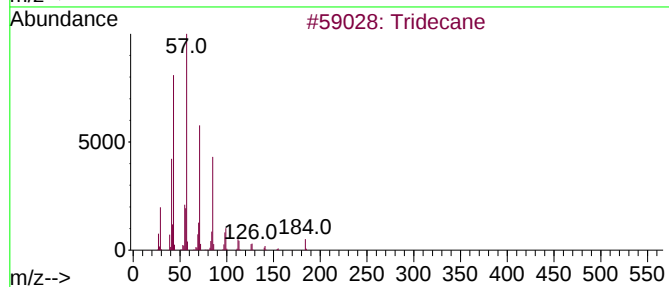
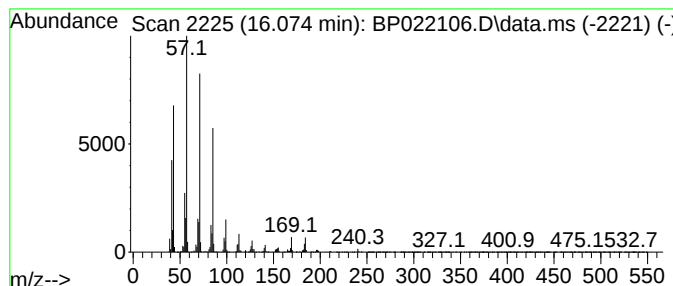
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 61 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.074	15.26 ng/ul	2181740	Phenanthrene-d10	17.127

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	90
2			Tridecane	184	C13H28	000629-50-5	90
3			Heptadecane	240	C17H36	000629-78-7	90
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	87
5			Nonadecane	268	C19H40	000629-92-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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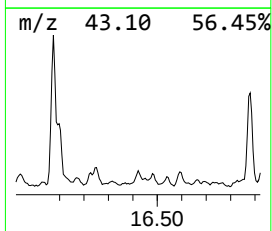
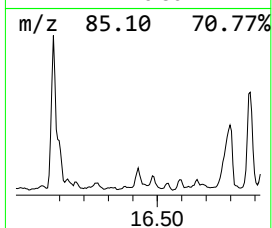
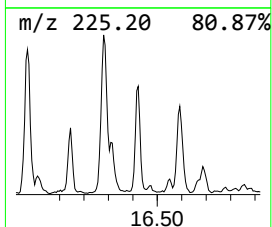
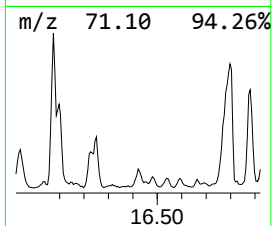
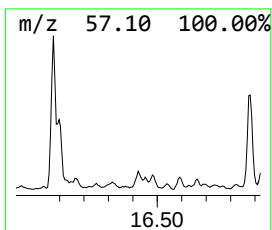
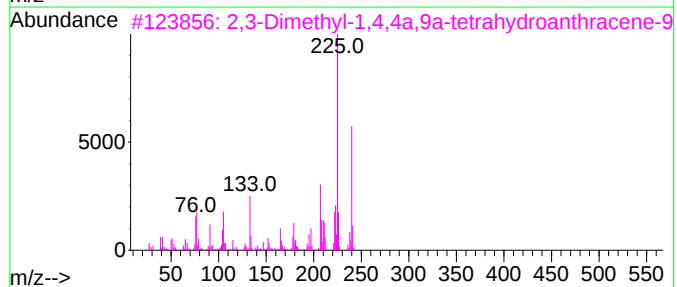
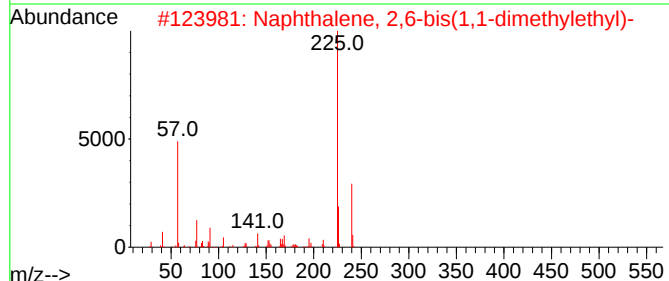
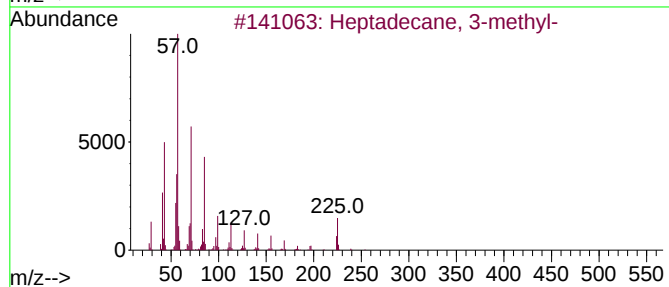
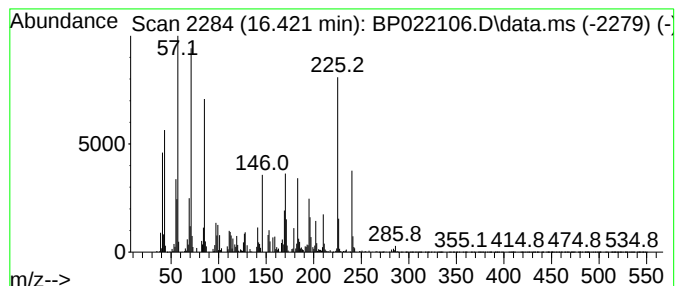
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 63 (DEL) Alkane: Straight-Chai... Concentration Rank 70

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.422	3.69 ng/ul	527664	Phenanthrene-d10	17.127

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 3-methyl-	254	C18H38	006418-44-6	35
2			Naphthalene, 2,6-bis(1,1-dimethy...	240	C18H24	003905-64-4	25
3			2,3-Dimethyl-1,4,4a,9a-tetrahydr...	240	C16H16O2	002670-23-7	25
4			4-(Dimethylamino)thiophenol, S-t...	225	C11H19NSSi	065257-43-4	22
5			Hexadecane, 5-butyl-	282	C20H42	006912-07-8	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC29DL

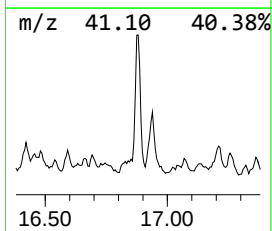
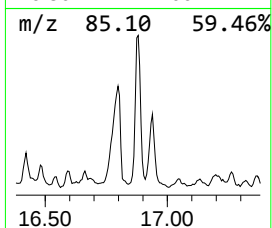
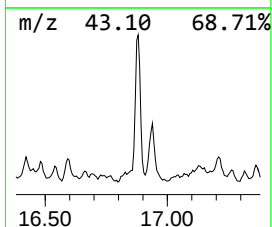
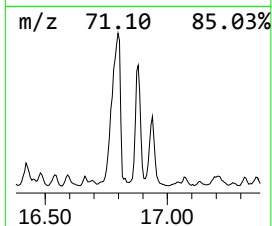
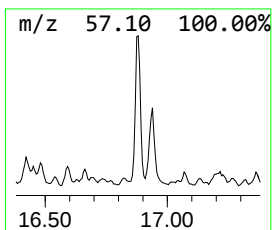
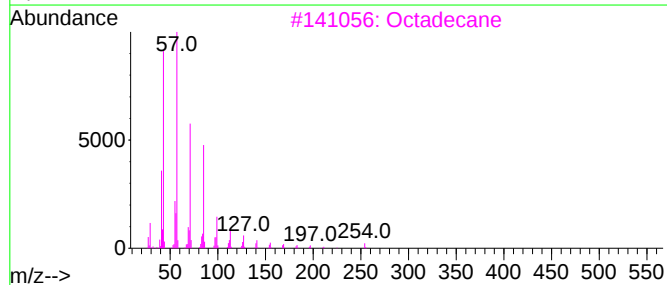
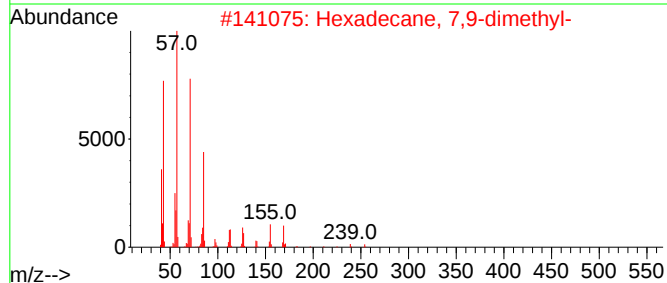
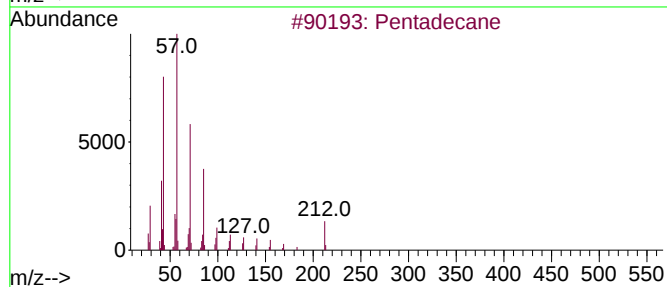
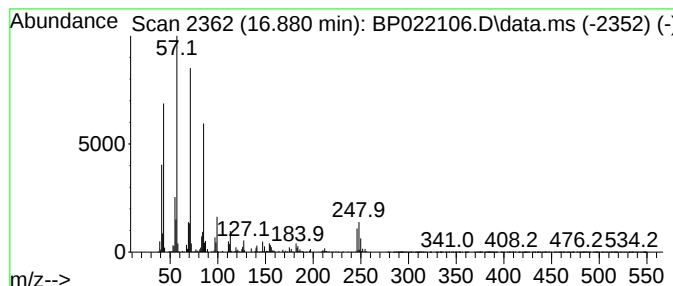
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 65 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.880	20.49 ng/ul	2929860	Phenanthrene-d10	17.127

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane	212	C15H32	000629-62-9	87
2			Hexadecane, 7,9-dimethyl-	254	C18H38	021164-95-4	86
3			Octadecane	254	C18H38	000593-45-3	80
4			Decane, 3-methyl-	156	C11H24	013151-34-3	64
5			Tridecane	184	C13H28	000629-50-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC29DL

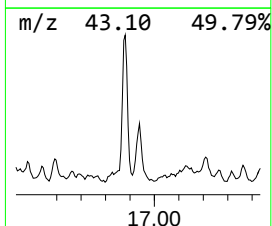
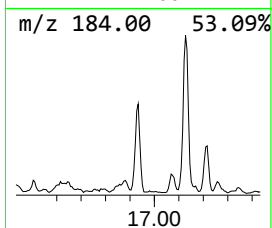
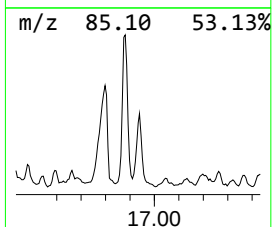
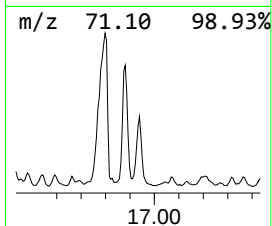
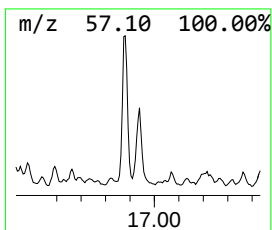
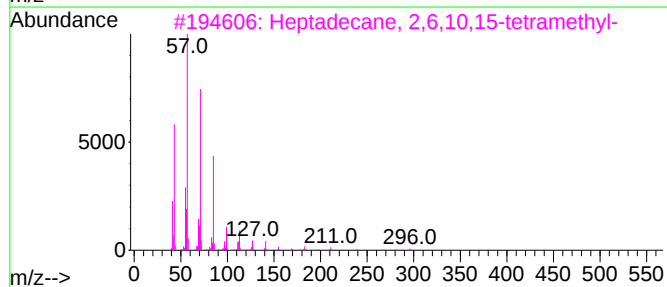
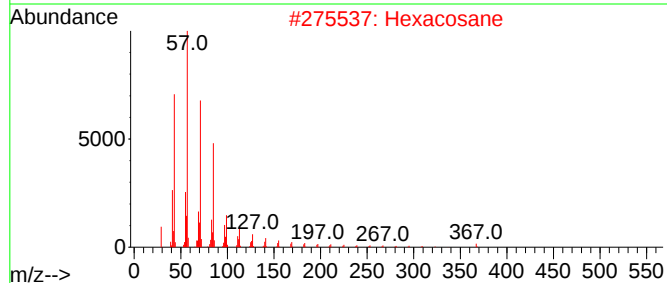
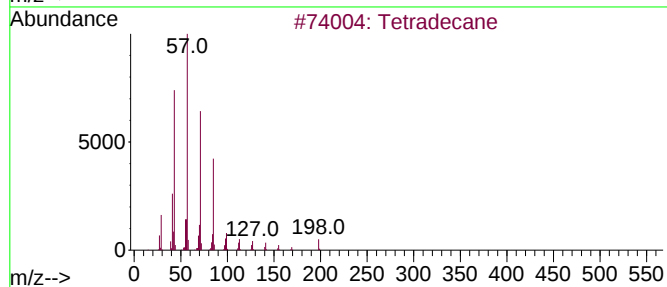
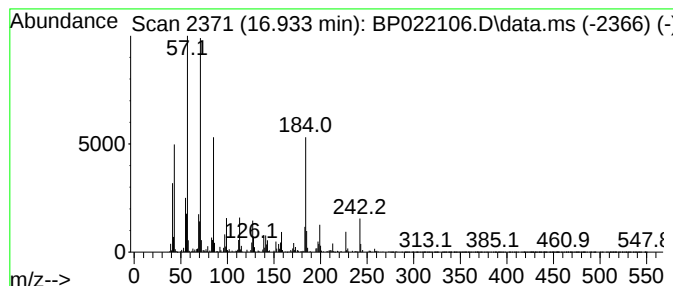
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 66 (DEL) Alkane: Straight-Chai... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.933	11.70 ng/ul	1673070	Phenanthrene-d10	17.127

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	55
2			Hexacosane	366	C26H54	000630-01-3	46
3			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	46
4			Tridecane	184	C13H28	000629-50-5	43
5			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC29DL

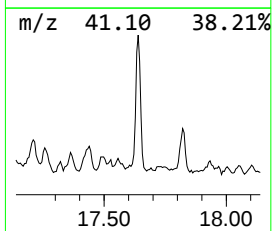
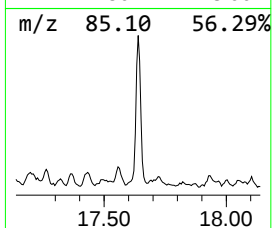
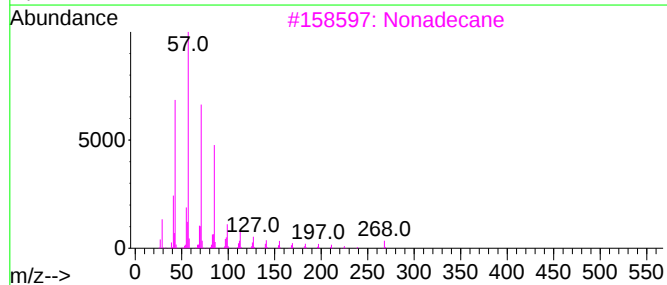
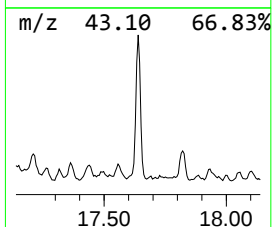
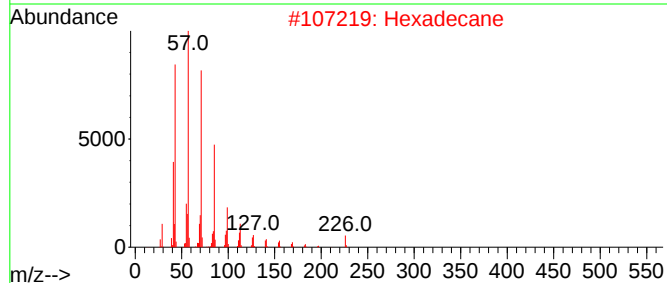
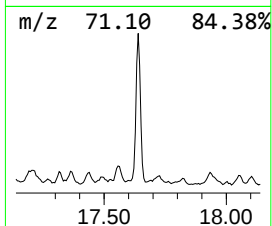
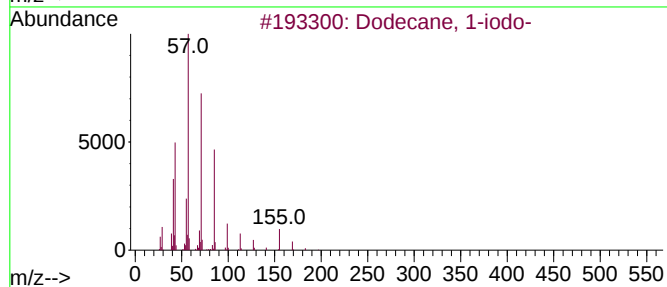
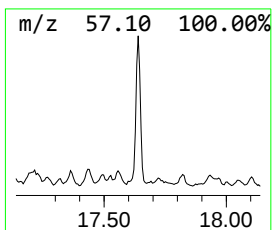
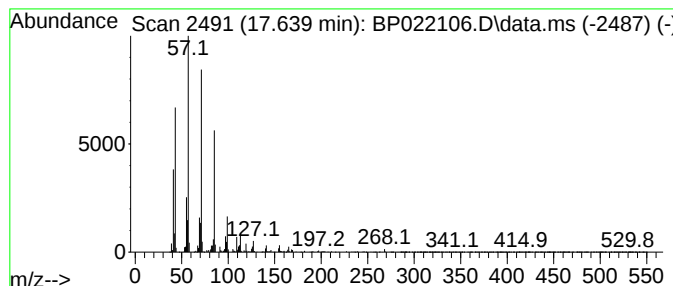
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 70 Dodecane, 1-iodo- Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.639	10.24 ng/ul	1464280	Phenanthrene-d10	17.127

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86
2			Hexadecane	226	C16H34	000544-76-3	86
3			Nonadecane	268	C19H40	000629-92-5	86
4			Tetracosane	338	C24H50	000646-31-1	86
5			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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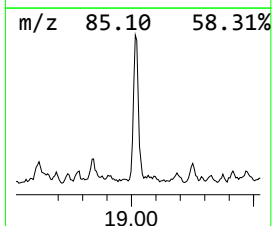
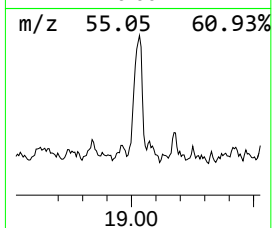
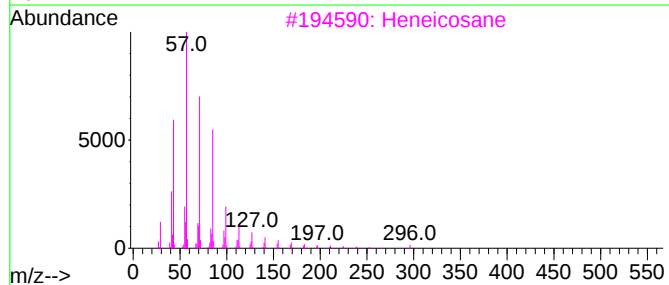
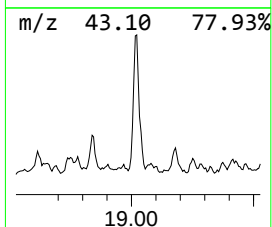
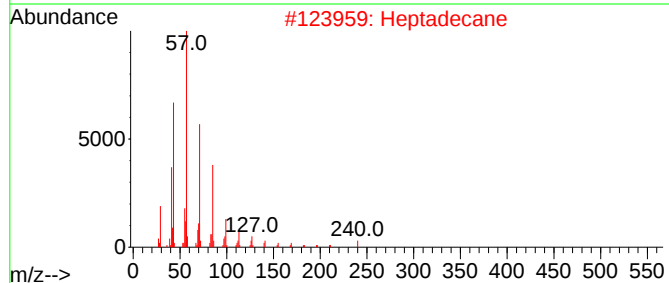
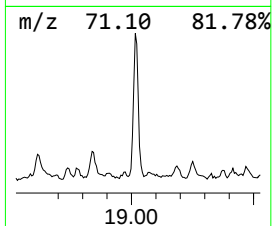
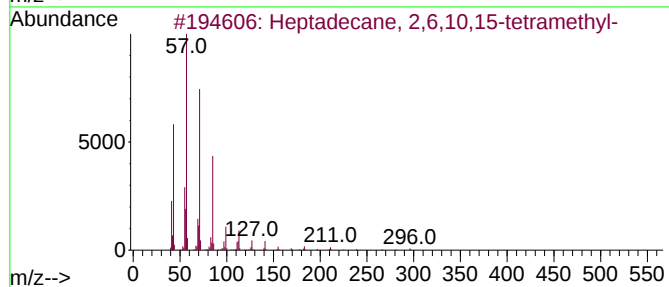
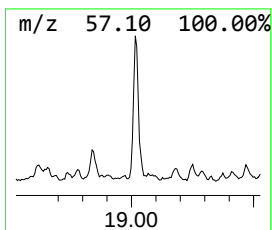
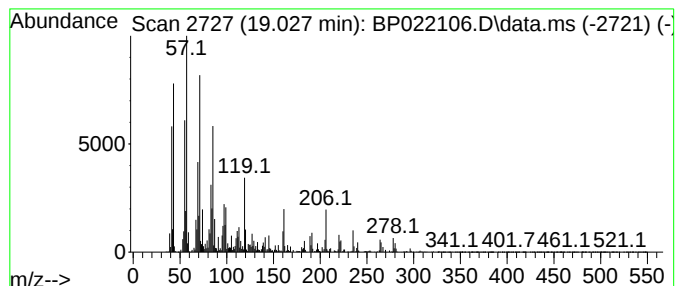
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 74 (DEL) Alkane: Straight-Chai... Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.027	11.34 ng/ul	1621700	Phenanthrene-d10	17.127

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	90
2			Heptadecane	240	C17H36	000629-78-7	87
3			Heneicosane	296	C21H44	000629-94-7	87
4			Heptadecane	240	C17H36	000629-78-7	70
5			Octadecane	254	C18H38	000593-45-3	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

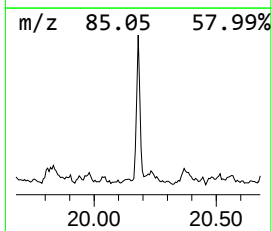
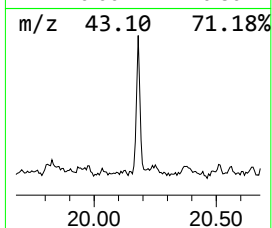
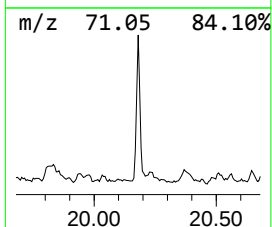
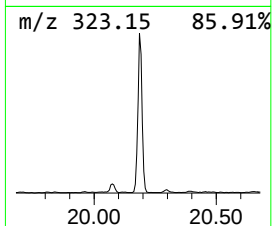
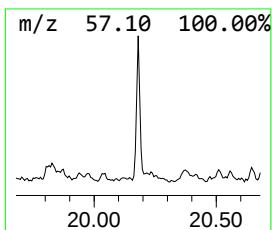
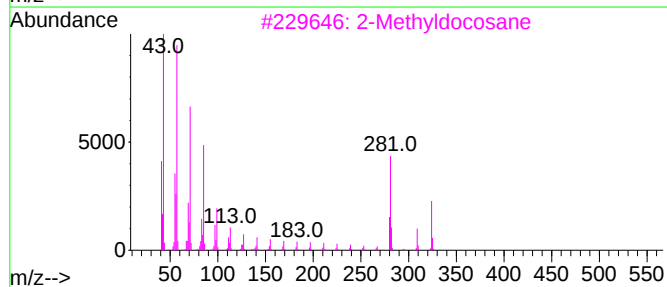
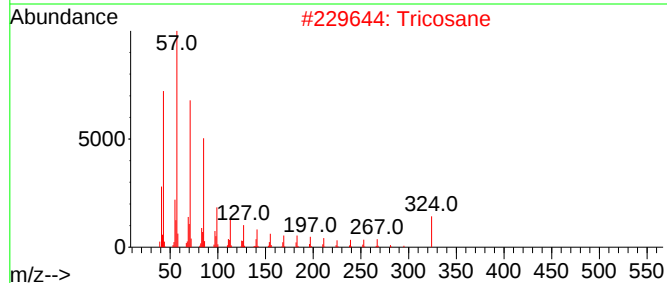
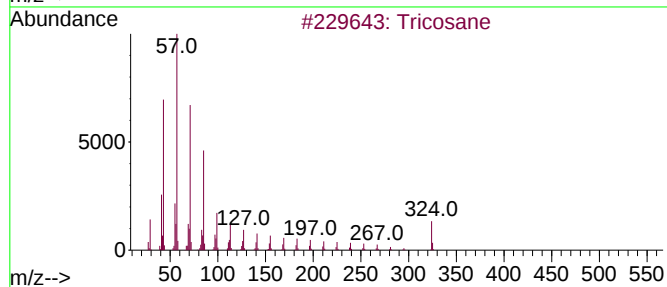
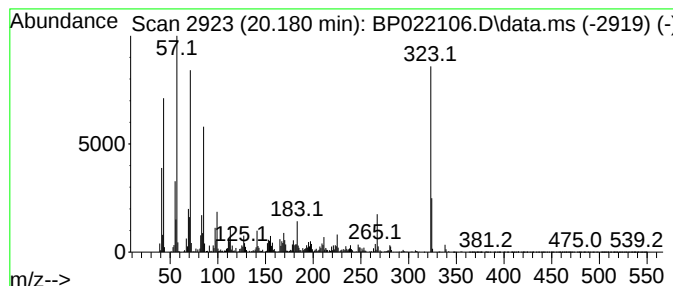
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 75 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.180	12.72 ng/ul	1273170	Chrysene-d12	21.545	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricosane	324	C23H48	000638-67-5	70
2	Tricosane	324	C23H48	000638-67-5	49
3	2-Methyldocosane	324	C23H48	001560-81-2	46
4	4-Methylheneicosane	310	C22H46	025117-29-7	43
5	5,5-Diethylheptadecane	296	C21H44	1010360-41-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC29DL

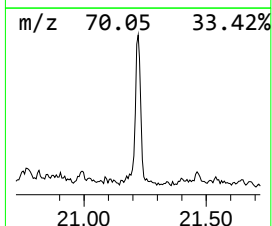
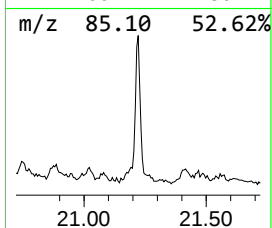
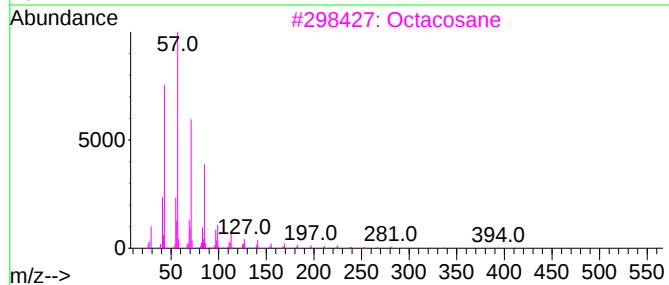
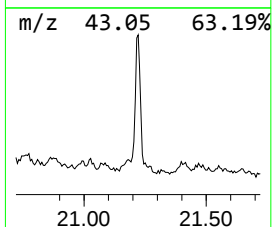
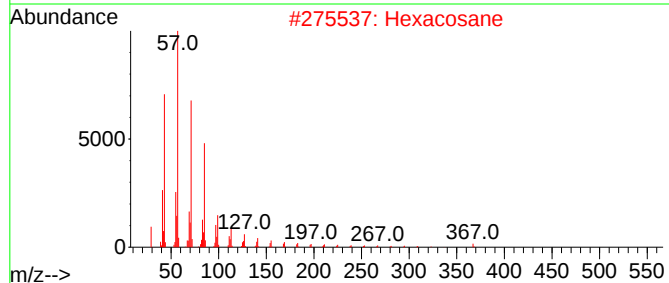
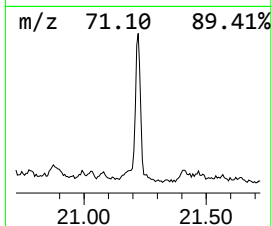
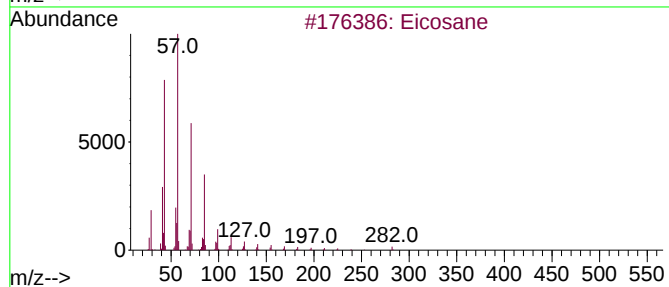
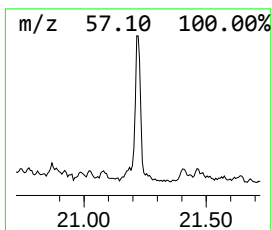
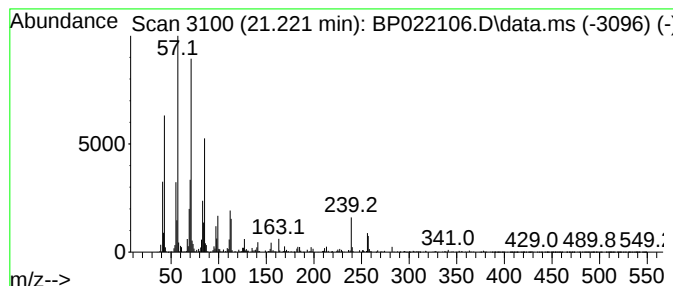
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 76 (DEL) Alkane: Straight-Chai... Concentration Rank 60

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.221	6.18 ng/ul	618270	Chrysene-d12	21.545

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	86
2		Hexacosane	366	C26H54	000630-01-3	81
3		Octacosane	394	C28H58	000630-02-4	68
4		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	64
5		Hexadecane	226	C16H34	000544-76-3	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022106.D
Acq On : 28 Sep 2024 15:46
Operator : RC/JU
Sample : P3910-10DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC29DL

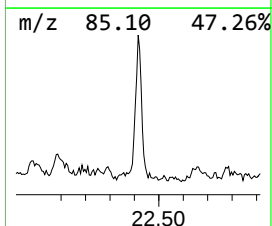
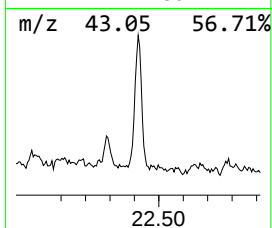
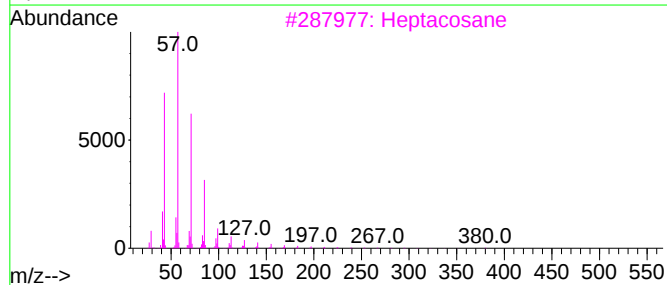
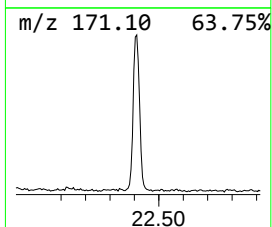
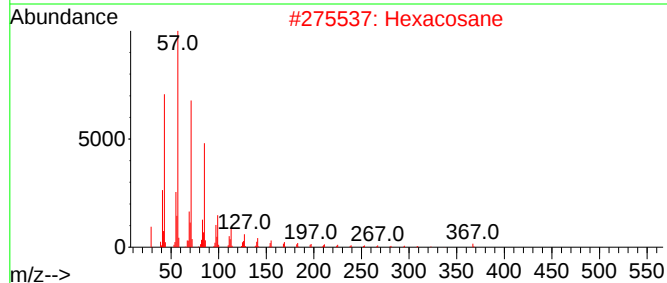
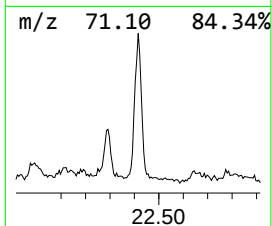
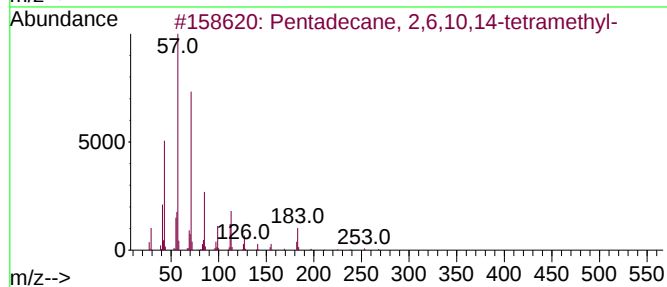
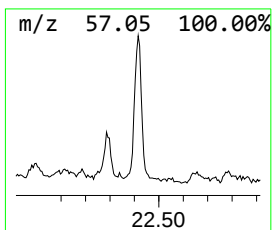
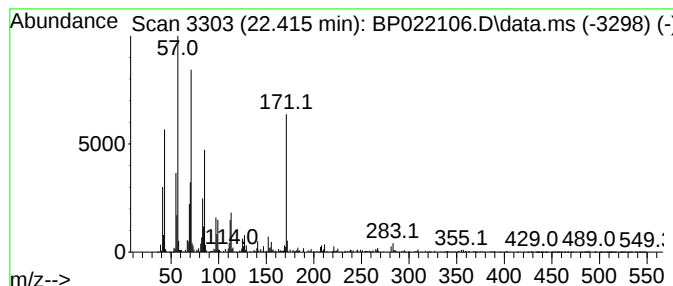
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 77 (DEL) Alkane: Straight-Chai... Concentration Rank 58

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.415	6.90 ng/ul	691023	Chrysene-d12	21.545

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	83
2		Hexacosane	366	C26H54	000630-01-3	68
3		Heptacosane	380	C27H56	000593-49-7	64
4		Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	64
5		Pentatriacontane	493	C35H72	000630-07-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
 Data File : BP022106.D
 Acq On : 28 Sep 2024 15:46
 Operator : RC/JU
 Sample : P3910-10DL 10X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DDC29DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	5.757	26.6	ng/ul	1815390	1	7.698	1366200	20.0
(DEL) Alkane: S...	6.287	7.0	ng/ul	480035	1	7.698	1366200	20.0
(DEL) Alkane: S...	6.710	16.7	ng/ul	1139930	1	7.698	1366200	20.0
1-Iodo-2-methyl...	6.769	11.7	ng/ul	800150	1	7.698	1366200	20.0
Benzene, 1-ethyl...	6.804	15.0	ng/ul	1022430	1	7.698	1366200	20.0
Benzene, 1-ethyl...	7.098	12.1	ng/ul	824139	1	7.698	1366200	20.0
2-Heptanone, 4,...	7.134	9.1	ng/ul	624488	1	7.698	1366200	20.0
(DEL) Alkane: S...	7.334	80.4	ng/ul	5489660	1	7.698	1366200	20.0
(DEL) Alkane: S...	7.622	8.4	ng/ul	575266	1	7.698	1366200	20.0
1-Hexanol, 2-et...	7.769	34.6	ng/ul	2365960	1	7.698	1366200	20.0
Benzene, 1,2,3-...	7.816	11.3	ng/ul	773555	1	7.698	1366200	20.0
Hexanal ethyl t...	7.916	10.9	ng/ul	741001	1	7.698	1366200	20.0
(DEL) Alkane: C...	7.987	7.1	ng/ul	481988	1	7.698	1366200	20.0
(DEL) Alkane: S...	8.228	12.4	ng/ul	845738	1	7.698	1366200	20.0
Benzene, 1,4-di...	8.328	16.8	ng/ul	1144660	1	7.698	1366200	20.0
(DEL) Alkane: S...	8.457	12.3	ng/ul	838997	1	7.698	1366200	20.0
Benzene, 1-meth...	8.492	10.7	ng/ul	729664	1	7.698	1366200	20.0
o-Cymene	8.687	10.4	ng/ul	709986	1	7.698	1366200	20.0
(DEL) Alkane: S...	8.910	59.9	ng/ul	4092970	1	7.698	1366200	20.0
unknown-01	9.034	13.5	ng/ul	920244	1	7.698	1366200	20.0
(DEL) Alkane: C...	9.604	3.7	ng/ul	832860	2	10.463	4498230	20.0
(DEL) Alkane: S...	9.751	2.5	ng/ul	573819	2	10.463	4498230	20.0
(DEL) Alkane: S...	9.892	4.5	ng/ul	1001900	2	10.463	4498230	20.0
(DEL) Alkane: C...	10.163	2.3	ng/ul	519123	2	10.463	4498230	20.0
(DEL) Alkane: S...	11.492	4.5	ng/ul	1010070	2	10.463	4498230	20.0
Benzene, 4-(2-b...	11.557	16.2	ng/ul	3641220	2	10.463	4498230	20.0
(DEL) Alkane: C...	11.928	4.8	ng/ul	1073400	2	10.463	4498230	20.0
(DEL) Alkane: S...	12.534	13.7	ng/ul	920632	3	14.316	1349190	20.0
Propanoic acid,...	12.722	12.1	ng/ul	817133	3	14.316	1349190	20.0
Butanoic acid, ...	13.057	14.4	ng/ul	970255	3	14.316	1349190	20.0
Butanoic acid, ...	13.269	23.7	ng/ul	1598230	3	14.316	1349190	20.0
Naphthalene, 2,...	13.486	20.6	ng/ul	1390670	3	14.316	1349190	20.0
Propanoic acid,...	13.575	24.8	ng/ul	1675920	3	14.316	1349190	20.0
Naphthalene, 1,...	13.633	9.4	ng/ul	634087	3	14.316	1349190	20.0
Bicyclo[4.1.0]h...	13.669	28.3	ng/ul	1911120	3	14.316	1349190	20.0
(DEL) Alkane: S...	13.804	9.6	ng/ul	649577	3	14.316	1349190	20.0
Propanoic acid,...	14.069	18.6	ng/ul	1258150	3	14.316	1349190	20.0
Oxalic acid, 2-...	14.227	119.0	ng/ul	8027580	3	14.316	1349190	20.0
unknown-02	14.404	15.7	ng/ul	1055650	3	14.316	1349190	20.0
4'-(1-Pyrrolyl)a...	14.463	70.5	ng/ul	4754240	3	14.316	1349190	20.0
Cyclohexanone, ...	14.710	26.3	ng/ul	1774300	3	14.316	1349190	20.0
4b,5,6,7,8,8a,9...	14.886	13.1	ng/ul	881643	3	14.316	1349190	20.0
Cyclopropane, 1...	15.092	33.6	ng/ul	2269180	3	14.316	1349190	20.0
unknown-03	15.163	11.3	ng/ul	765031	3	14.316	1349190	20.0
(DEL) Alkane: S...	15.186	28.2	ng/ul	1902440	3	14.316	1349190	20.0
3-Methyl-4-(met...	15.598	14.2	ng/ul	959929	3	14.316	1349190	20.0
unknown-04	15.698	12.1	ng/ul	817679	3	14.316	1349190	20.0
(DEL) Alkane: S...	16.074	15.3	ng/ul	2181740	4	17.127	2859420	20.0
(DEL) Alkane: S...	16.422	3.7	ng/ul	527664	4	17.127	2859420	20.0
(DEL) Alkane: S...	16.880	20.5	ng/ul	2929860	4	17.127	2859420	20.0
(DEL) Alkane: S...	16.933	11.7	ng/ul	1673070	4	17.127	2859420	20.0
Dodecane, 1-iodo-	17.639	10.2	ng/ul	1464280	4	17.127	2859420	20.0

(DEL) Alkane: S...	19.027	11.3	ng/ul	1621700	4	17.127	2859420	20.0
(DEL) Alkane: S...	20.180	12.7	ng/ul	1273170	5	21.545	2002060	20.0
(DEL) Alkane: S...	21.221	6.2	ng/ul	618270	5	21.545	2002060	20.0
(DEL) Alkane: S...	22.415	6.9	ng/ul	691023	5	21.545	2002060	20.0

Instrument :
BNA_P
ClientSampleId :
DDC29DL