

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
 Data File : BP022100.D
 Acq On : 28 Sep 2024 11:42
 Operator : RC/JU
 Sample : P3910-02ME
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DDC75ME

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: BP022100.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.758	464	471	479	rVB2	696798	1088829	2.05%	0.631%
2	6.369	566	575	578	rVV3	225342	542179	1.02%	0.314%
3	6.711	623	633	639	rBV3	521185	1234256	2.33%	0.715%
4	6.769	639	643	646	rVV	547066	787313	1.49%	0.456%
5	6.811	646	650	655	rVV	948735	1464566	2.76%	0.849%
6	6.875	655	661	667	rVV4	1038463	2159761	4.08%	1.252%
7	6.940	667	672	678	rVB	487142	713502	1.35%	0.414%
8	7.105	694	700	703	rVV	547714	875462	1.65%	0.507%
9	7.140	703	706	714	rVV	650248	1095934	2.07%	0.635%
10	7.246	719	724	734	rVV2	285806	650867	1.23%	0.377%
11	7.340	734	740	748	rVB2	2988443	5825622	10.99%	3.377%
12	7.628	778	789	793	rBV5	189182	561012	1.06%	0.325%
13	7.687	793	799	805	rVB2	660279	1280032	2.42%	0.742%
14	7.769	805	813	817	rBV3	387273	767510	1.45%	0.445%
15	7.816	817	821	829	rVB3	459979	838166	1.58%	0.486%
16	7.993	847	851	857	rVB3	289878	532355	1.00%	0.309%
17	8.128	868	874	878	rBV	1322925	2129731	4.02%	1.234%
18	8.175	878	882	887	rBV2	276750	466535	0.88%	0.270%
19	8.234	887	892	897	rVB2	517631	1030397	1.94%	0.597%
20	8.334	904	909	911	rBV2	774997	1218220	2.30%	0.706%
21	8.393	916	919	925	rVB2	361752	632655	1.19%	0.367%
22	8.458	925	930	933	rBV	646055	915367	1.73%	0.531%
23	8.493	933	936	944	rVB	475180	761872	1.44%	0.442%
24	8.640	956	961	965	rBV	404892	620905	1.17%	0.360%
25	8.687	965	969	977	rVB2	549444	994802	1.88%	0.577%
26	8.793	977	987	992	rBV2	647712	1393842	2.63%	0.808%
27	8.916	999	1008	1014	rVB	2576696	4059682	7.66%	2.353%
28	9.034	1018	1028	1035	rVB8	345422	939770	1.77%	0.545%
29	9.116	1035	1042	1049	rBV4	237750	736175	1.39%	0.427%
30	9.363	1080	1084	1094	rVB2	259333	602424	1.14%	0.349%
31	9.469	1094	1102	1108	rVB4	204337	492219	0.93%	0.285%
32	9.558	1109	1117	1121	rBV3	401877	856294	1.62%	0.496%
33	9.610	1121	1126	1130	rBV4	289710	496027	0.94%	0.288%
34	9.752	1147	1150	1155	rVB	360498	511533	0.97%	0.296%
35	9.816	1155	1161	1165	rBV3	273597	445190	0.84%	0.258%
36	10.087	1204	1207	1214	rVB2	422494	681735	1.29%	0.395%

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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

37	10.158	1214	1219	1230	rVB4	271217	536874	1.01%	0.311%
38	10.458	1259	1270	1276	rBV2	3779769	7431055	14.02%	4.307%
39	10.581	1285	1291	1304	rVB5	1094834	2561403	4.83%	1.485%
40	10.840	1329	1335	1345	rBV	857087	1467362	2.77%	0.851%
41	10.969	1348	1357	1363	rBV	609614	1164772	2.20%	0.675%
42	11.369	1417	1425	1431	rBV6	276087	795126	1.50%	0.461%
43	11.487	1439	1445	1450	rBV	636918	1033936	1.95%	0.599%
44	11.722	1475	1485	1493	rBV7	380534	851214	1.61%	0.493%
45	11.881	1506	1512	1516	rBV	824704	1312826	2.48%	0.761%
46	11.916	1516	1518	1523	rVB	371392	460097	0.87%	0.267%
47	12.128	1547	1554	1560	rVB3	740286	1505110	2.84%	0.872%
48	12.293	1577	1582	1587	rBV3	728539	1471115	2.78%	0.853%
49	12.557	1623	1627	1635	rVV4	258018	561272	1.06%	0.325%
50	12.646	1636	1642	1647	rVV6	193649	466661	0.88%	0.270%
51	12.704	1647	1652	1662	rVV7	196271	531867	1.00%	0.308%
52	12.881	1679	1682	1686	rVV	348450	498566	0.94%	0.289%
53	13.134	1720	1725	1732	rVV	978082	1621927	3.06%	0.940%
54	13.269	1745	1748	1754	rVV2	184658	440130	0.83%	0.255%
55	13.357	1755	1763	1770	rVB4	477879	1173805	2.21%	0.680%
56	13.487	1775	1785	1790	rBV4	631409	1647432	3.11%	0.955%
57	13.628	1803	1809	1813	rVV2	611300	1244525	2.35%	0.721%
58	13.675	1814	1817	1820	rVV2	434895	723329	1.36%	0.419%
59	13.716	1820	1824	1829	rVB	751292	1115920	2.11%	0.647%
60	13.799	1829	1838	1842	rBV4	274083	638868	1.21%	0.370%
61	13.863	1842	1849	1853	rBV4	384103	694931	1.31%	0.403%
62	13.999	1869	1872	1875	rVV	648671	906662	1.71%	0.526%
63	14.040	1875	1879	1884	rVB2	490273	717230	1.35%	0.416%
64	14.222	1906	1910	1914	rBV	1102174	1325485	2.50%	0.768%
65	14.310	1920	1925	1931	rVV	942186	1517051	2.86%	0.879%
66	14.387	1932	1938	1945	rVB5	340795	655087	1.24%	0.380%
67	14.534	1957	1963	1975	rBV6	577212	1573047	2.97%	0.912%
68	14.716	1990	1994	2000	rVB2	453775	807844	1.52%	0.468%
69	14.793	2000	2007	2012	rBV2	353872	624233	1.18%	0.362%
70	14.863	2013	2019	2023	rBV5	335455	684525	1.29%	0.397%
71	14.916	2023	2028	2029	rBV3	319306	433626	0.82%	0.251%
72	14.951	2029	2034	2039	rVB	3557641	4617555	8.71%	2.676%
73	15.187	2066	2074	2079	rVB2	1048078	1530494	2.89%	0.887%
74	15.316	2092	2096	2101	rVB	973407	1321651	2.49%	0.766%
75	15.434	2109	2116	2120	rBV5	454524	947810	1.79%	0.549%
76	15.604	2139	2145	2153	rVV3	457489	972861	1.84%	0.564%

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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 >
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77	15.693	2157	2160	2165	rVB3	346750	533998	1.01%	0.310%
78	16.063	2218	2223	2227	rBV	1095382	1682062	3.17%	0.975%
79	16.775	2334	2344	2354	rBV	28422781	52995291	100.00%	30.718%
80	16.881	2358	2362	2366	rVV	1013393	1267834	2.39%	0.735%
81	16.934	2366	2371	2380	rVB	589620	943159	1.78%	0.547%
82	17.110	2396	2401	2405	rBV	1152426	1586794	2.99%	0.920%
83	17.216	2413	2419	2425	rBV2	1002102	1776044	3.35%	1.029%
84	17.645	2487	2492	2496	rBV	1020085	1271229	2.40%	0.737%
85	17.710	2497	2503	2507	rBV	317345	518145	0.98%	0.300%
86	17.816	2516	2521	2525	rBV	1642652	2011499	3.80%	1.166%
87	18.051	2558	2561	2567	rVB2	338564	648015	1.22%	0.376%
88	18.363	2610	2614	2619	rBV	981179	1117612	2.11%	0.648%
89	19.022	2721	2726	2727	rBV2	906511	1132024	2.14%	0.656%
90	19.045	2727	2730	2734	rVB	1274716	1546127	2.92%	0.896%
91	19.181	2748	2753	2760	rVB2	708976	1025331	1.93%	0.594%
92	19.586	2817	2822	2826	rBV	1579203	2108847	3.98%	1.222%
93	19.628	2826	2829	2835	rVB2	599721	708192	1.34%	0.410%
94	20.186	2919	2924	2931	rBV	492160	693249	1.31%	0.402%
95	20.698	3008	3011	3019	rBV	456828	553054	1.04%	0.321%
96	21.222	3095	3100	3111	rVB	936413	1410160	2.66%	0.817%
97	21.539	3149	3154	3164	rVB2	1254799	2033585	3.84%	1.179%
98	21.786	3192	3196	3203	rVB	305996	520725	0.98%	0.302%
99	22.422	3298	3304	3312	rVB	567411	1092441	2.06%	0.633%
100	24.816	3703	3711	3723	rVB	943458	2360563	4.45%	1.368%

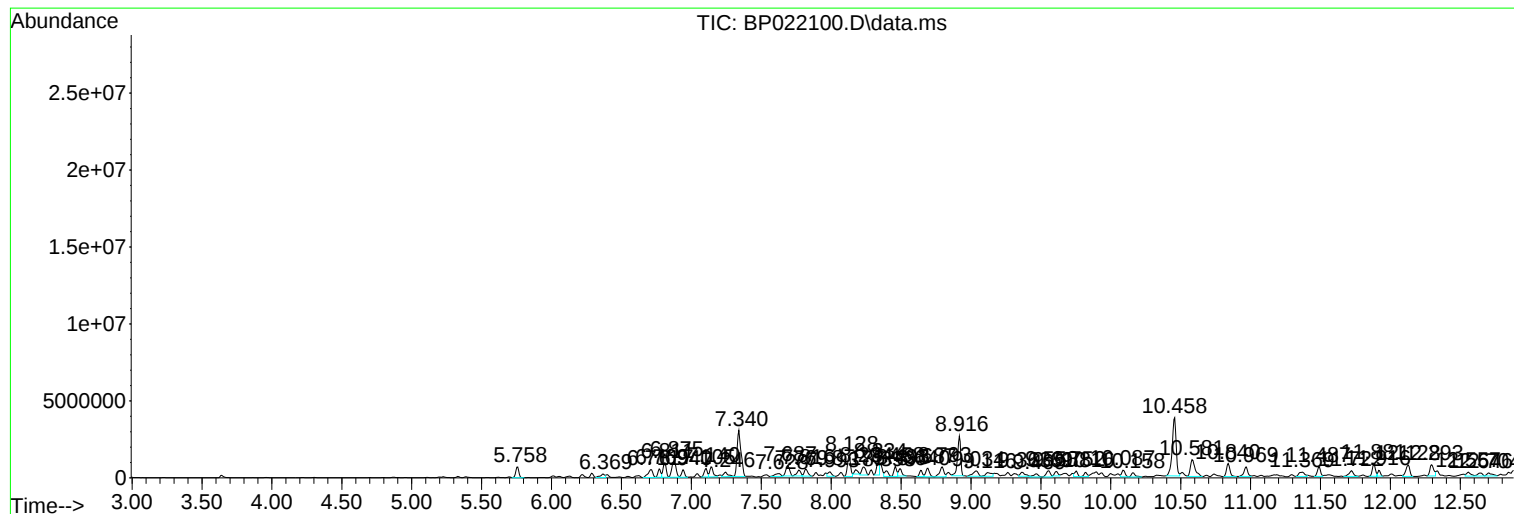
Sum of corrected areas: 172524005

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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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ALS Vial : 6 Sample Multiplier: 1

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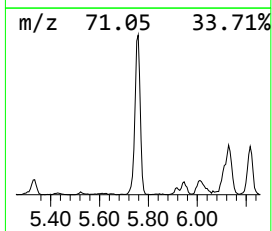
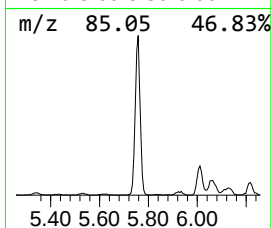
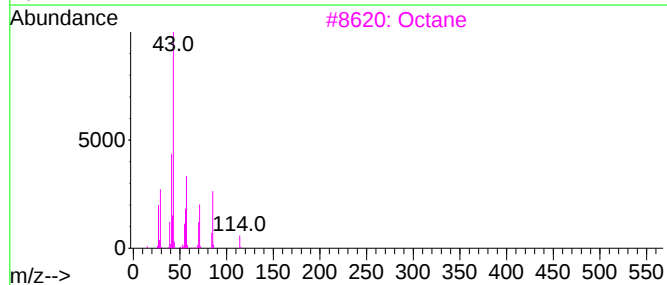
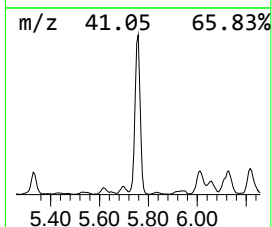
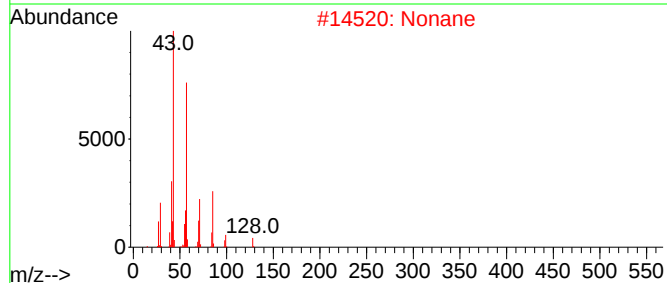
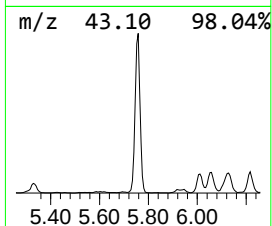
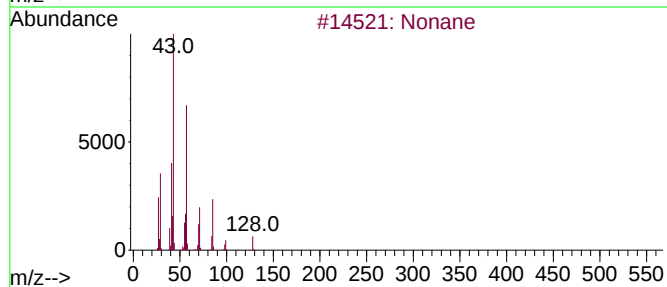
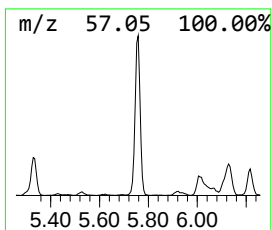
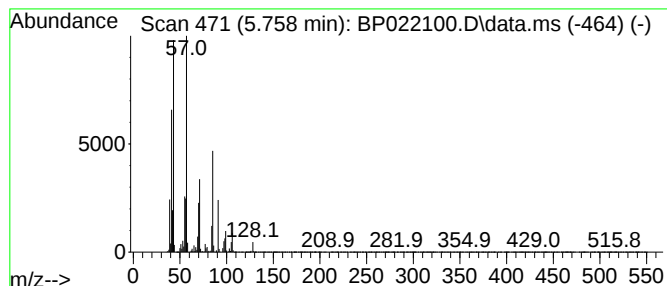
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.758	17.01 ng/ul	1088830	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	90
2			Nonane	128	C9H20	000111-84-2	68
3			Octane	114	C8H18	000111-65-9	59
4			Octane	114	C8H18	000111-65-9	50
5			Oxalic acid, isohexyl pentyl ester	244	C13H24O4	1000309-32-8	50



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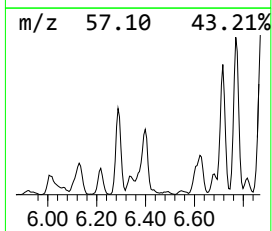
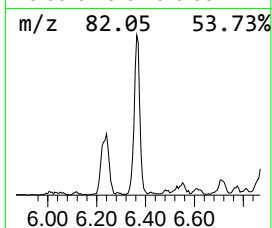
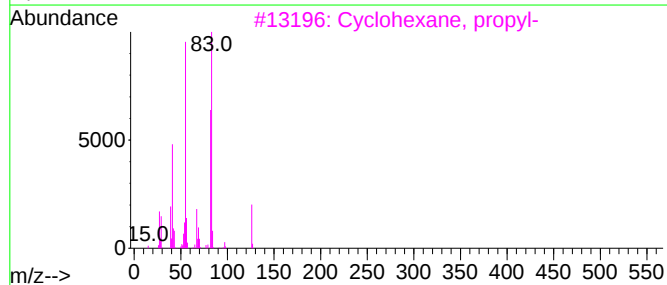
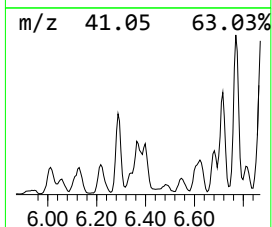
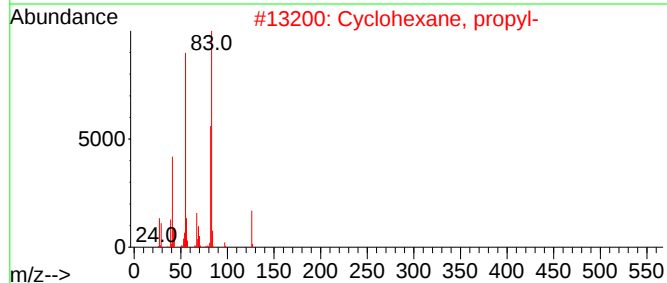
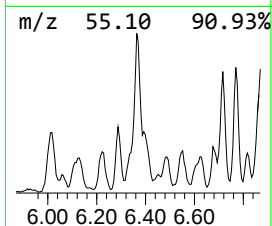
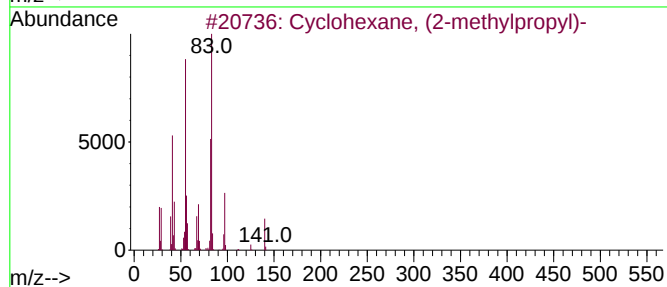
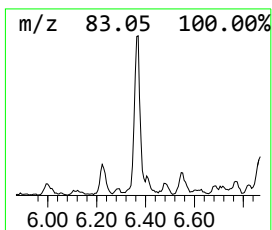
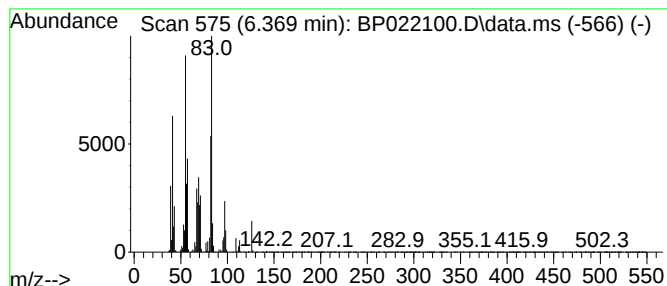
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Cyclic6.369 Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.369	8.47 ng/ul	542179	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	50
2			Cyclohexane, propyl-	126	C9H18	001678-92-8	49
3			Cyclohexane, propyl-	126	C9H18	001678-92-8	49
4			Undec-10-ynoic acid, dodecyl ester	350	C23H42O2	1010406-16-5	49
5			Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	43



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Misc :
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Instrument :
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ClientSampleId :
DDC75ME

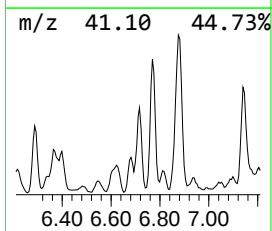
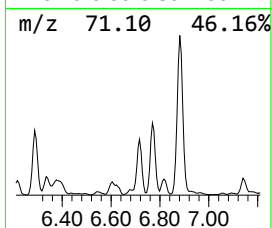
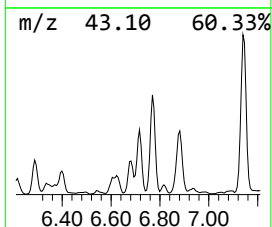
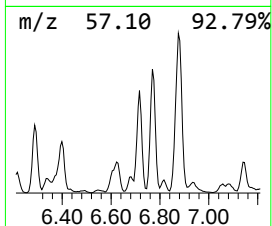
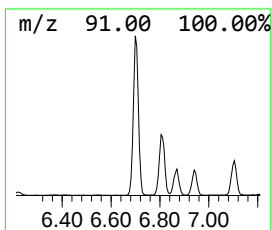
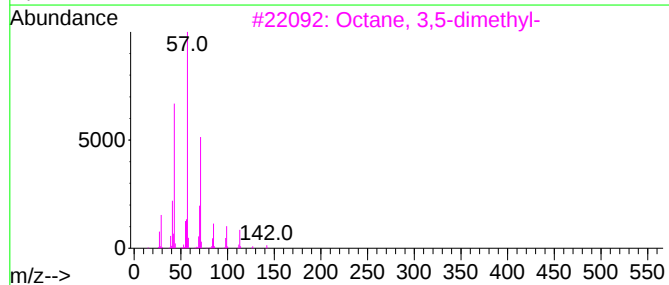
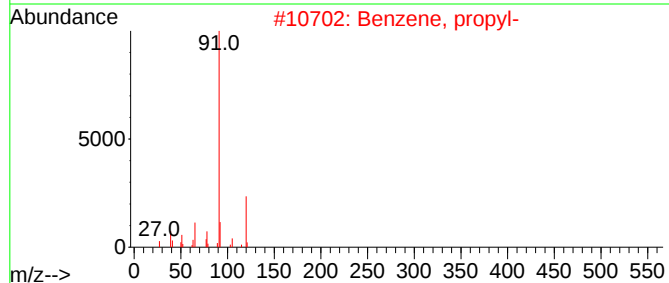
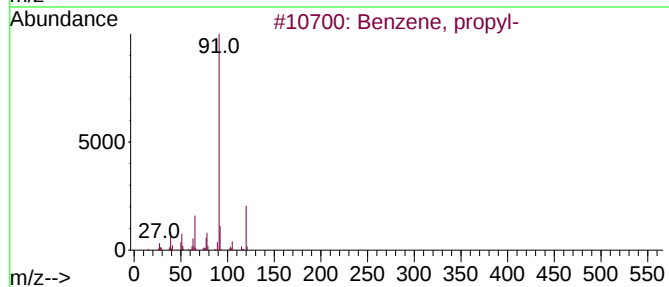
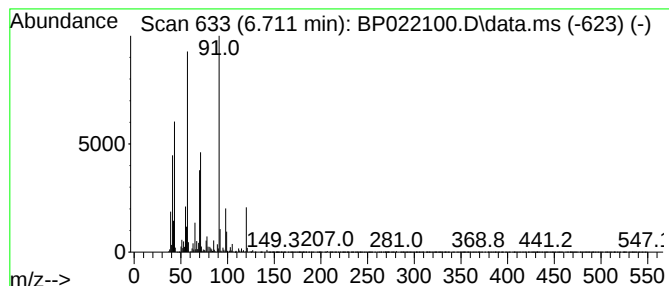
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, propyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.711	19.28 ng/ul	1234260	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	64
2			Benzene, propyl-	120	C9H12	000103-65-1	60
3			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	45
4			Nonane, 4-methyl-	142	C10H22	017301-94-9	43
5			Benzene, propyl-	120	C9H12	000103-65-1	35



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Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
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Instrument :
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ClientSampleId :
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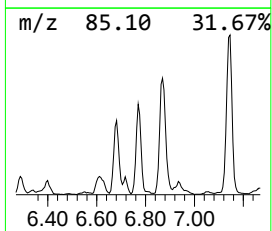
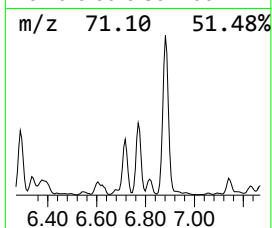
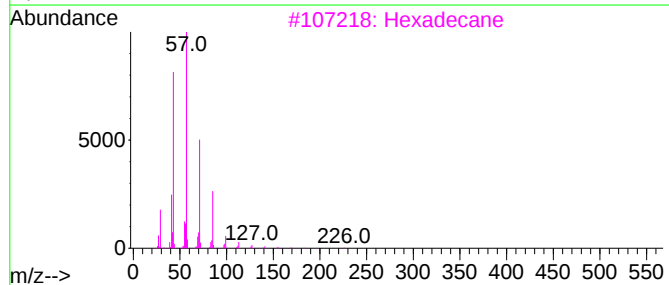
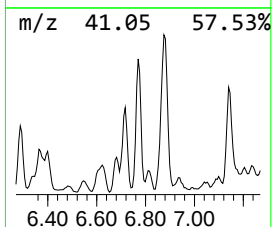
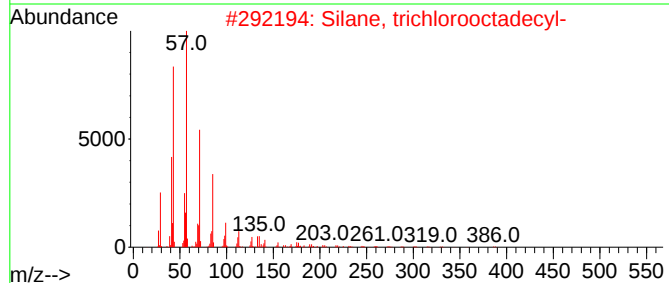
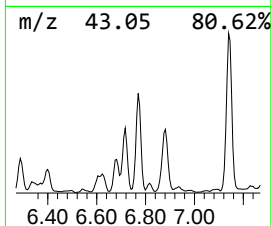
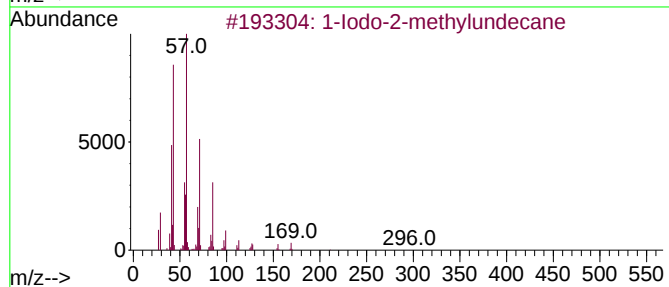
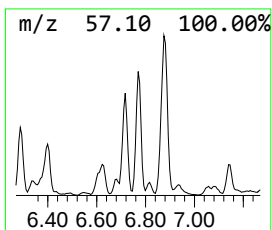
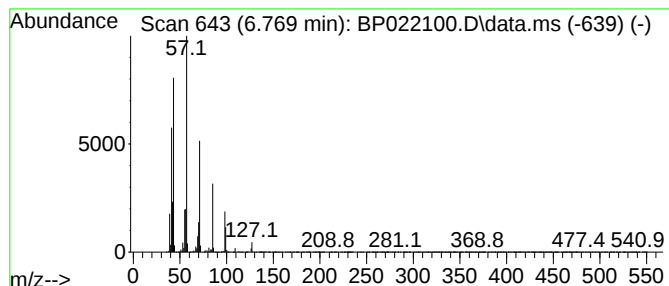
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Iodo-2-methylundecane Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.769	12.30 ng/ul	787313	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	78
2			Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	78
3			Hexadecane	226	C16H34	000544-76-3	64
4			Undecane, 2-methyl-	170	C12H26	007045-71-8	64
5			Undecane, 3,7-dimethyl-	184	C13H28	017301-29-0	53



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Acq On : 28 Sep 2024 11:42
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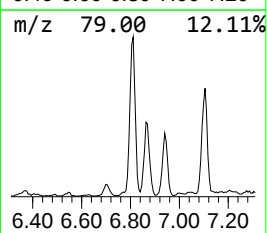
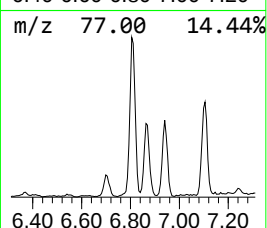
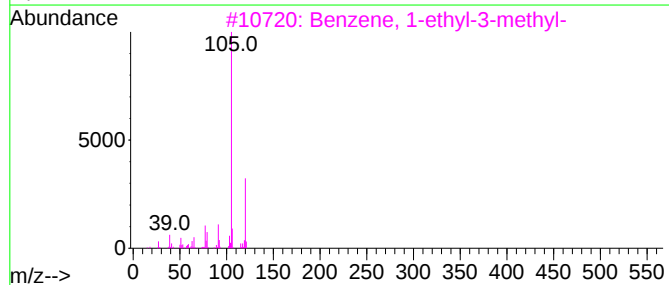
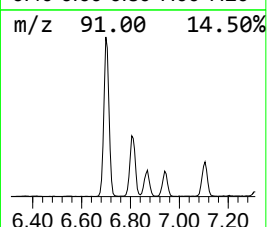
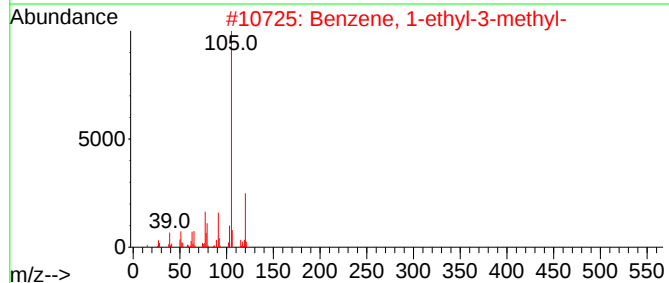
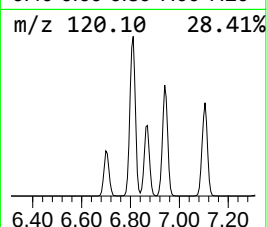
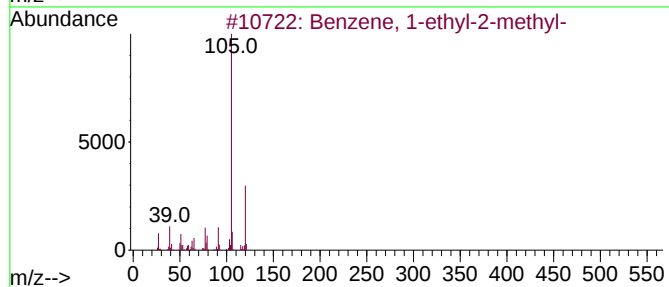
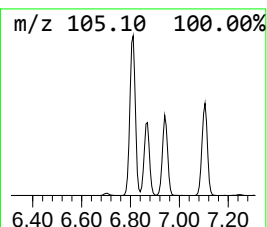
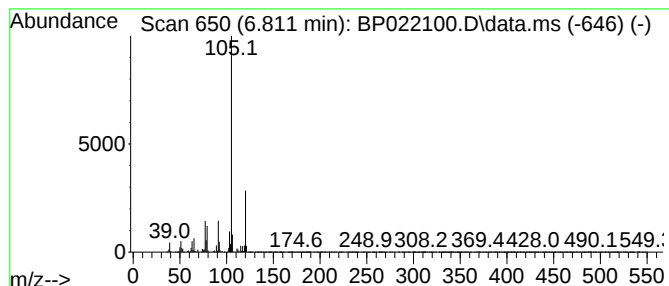
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.811	22.88 ng/ul	1464570	1,4-Dichlorobenzene-d4	7.699

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	95
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
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Acq On : 28 Sep 2024 11:42
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Sample : P3910-02ME
Misc :
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Instrument :
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ClientSampleId :
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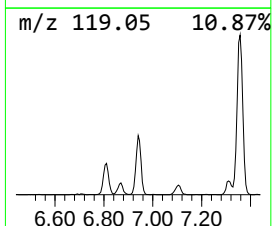
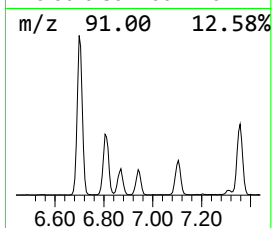
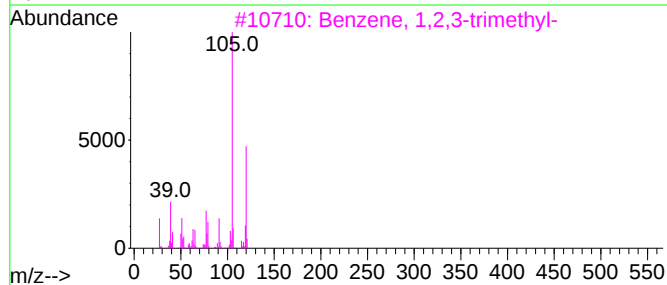
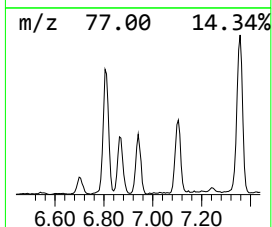
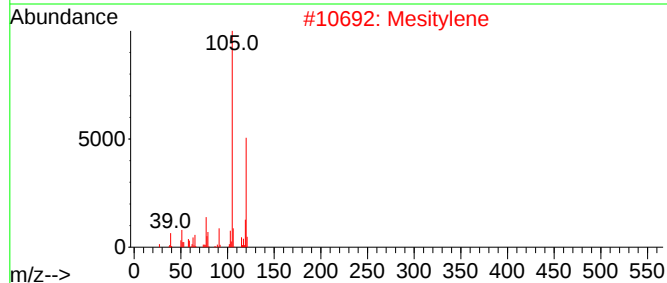
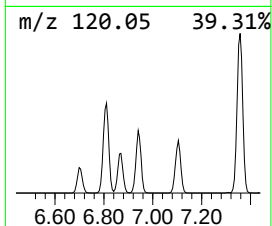
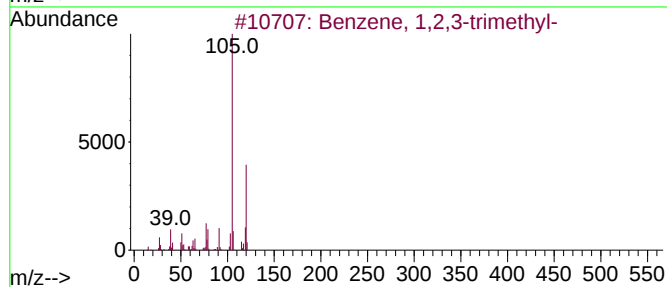
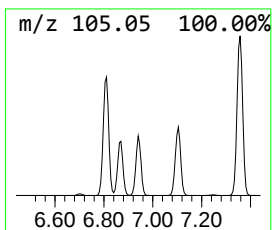
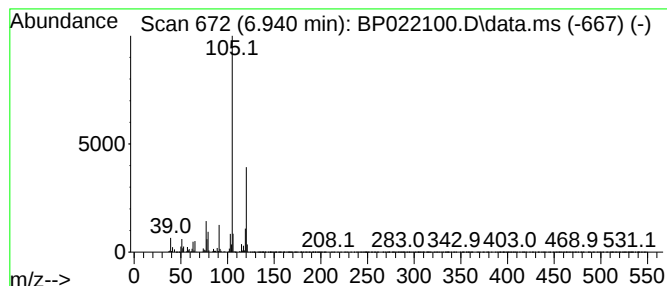
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.940	11.15 ng/ul	713502	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2			Mesitylene	120	C9H12	000108-67-8	95
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
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Sample : P3910-02ME
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Instrument :
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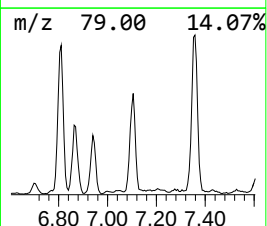
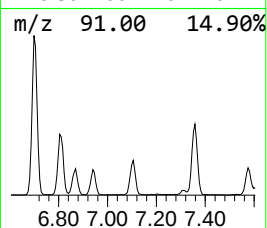
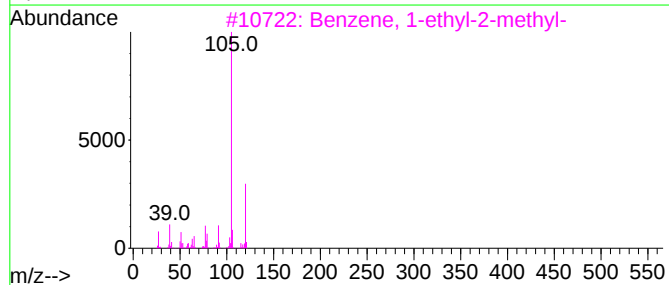
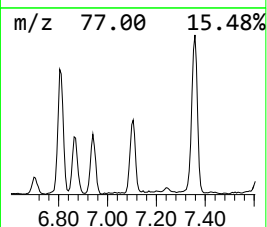
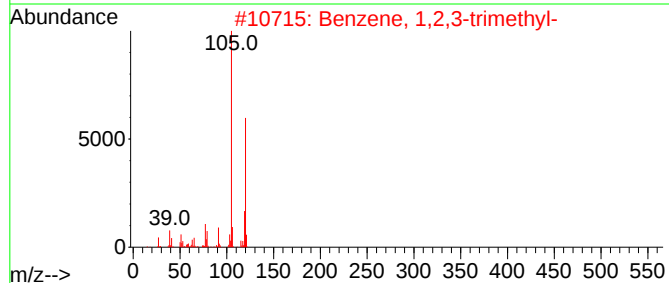
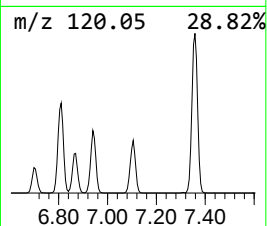
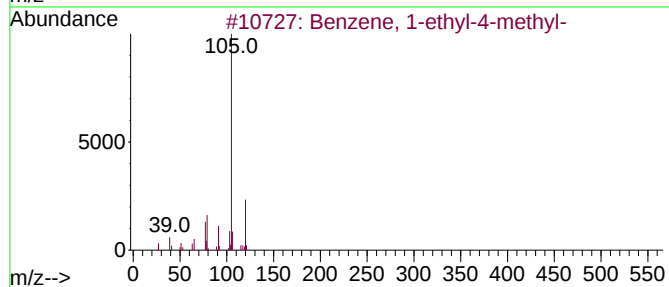
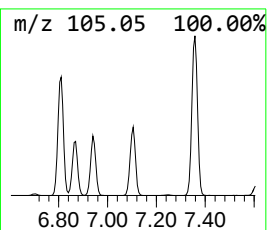
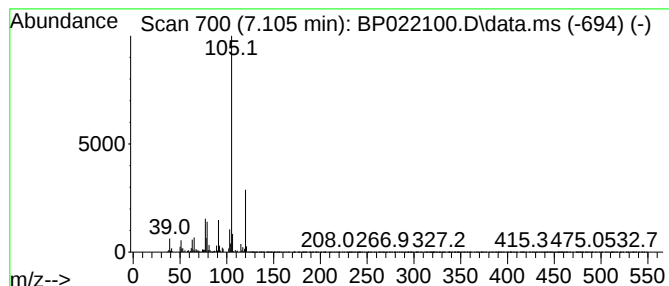
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1-ethyl-4-methyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.105	13.68 ng/ul	875462	1,4-Dichlorobenzene-d4	7.699

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
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Misc :
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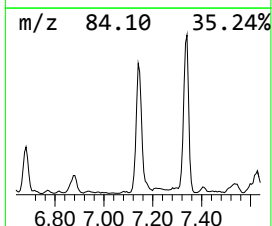
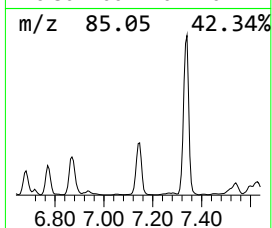
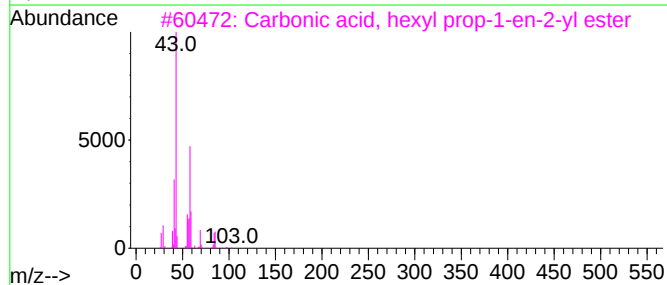
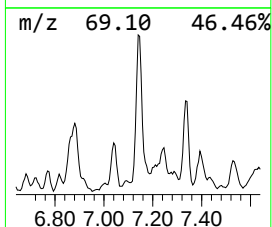
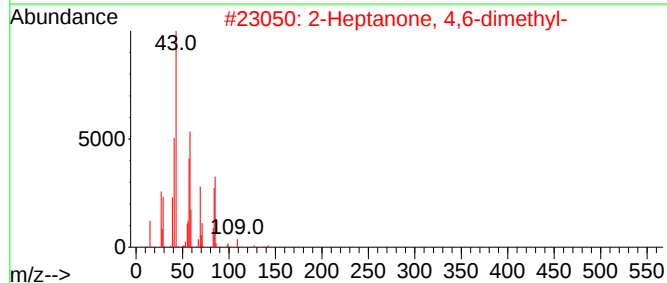
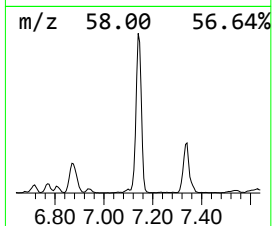
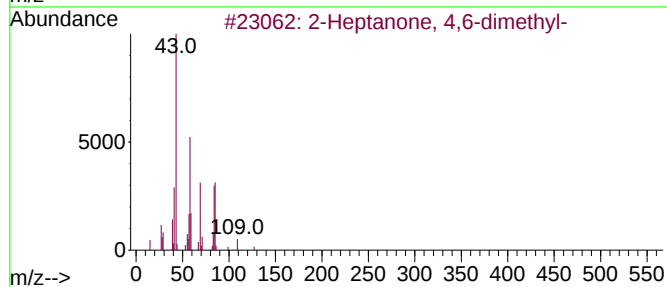
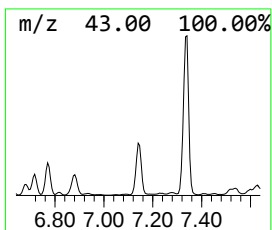
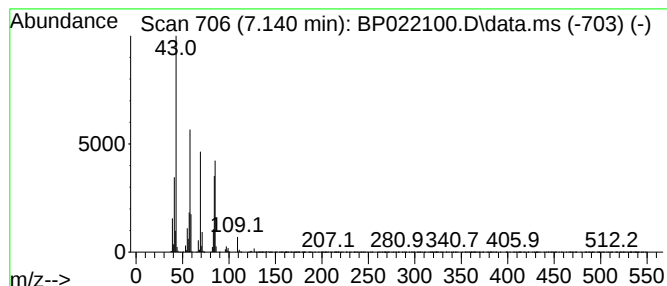
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2-Heptanone, 4,6-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.140	17.12 ng/ul	1095930	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	90		
2	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	64		
3	Carbonic acid, hexyl prop-1-en-2...	186	C10H18O3	1000382-54-2	59		
4	2-Nonanone, 9-[(tetrahydro-2H-py...	242	C14H26O3	054699-41-1	45		
5	Butanoic acid, 3-oxo-, 2-propeny...	142	C7H10O3	001118-84-9	45		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
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Misc :
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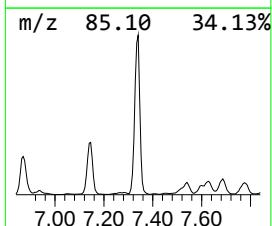
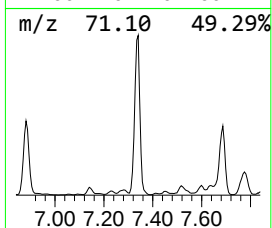
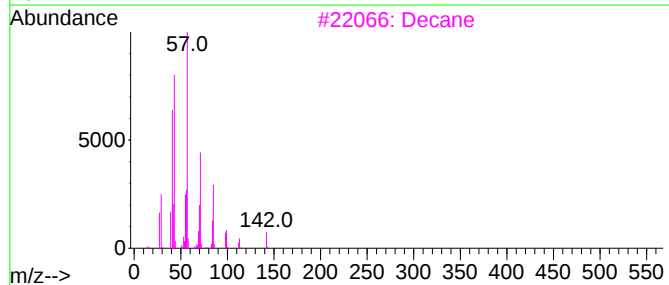
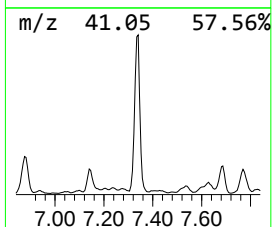
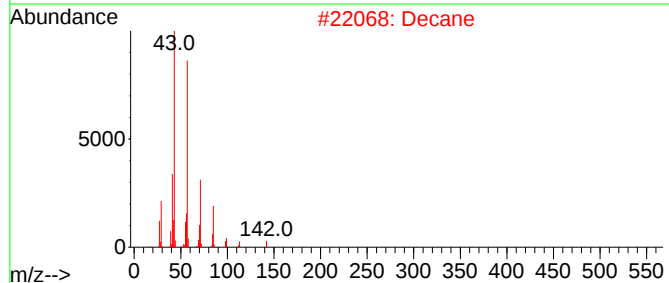
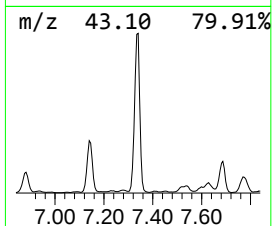
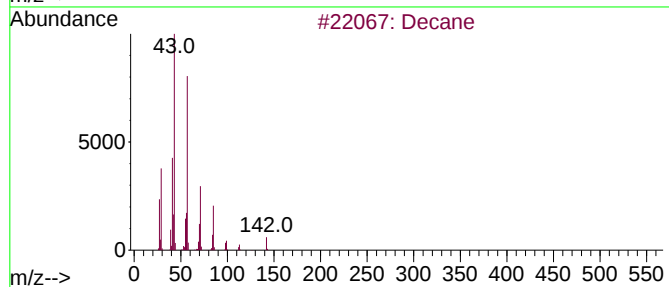
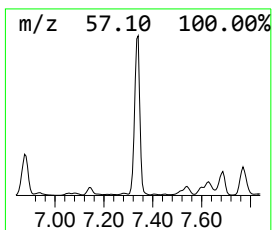
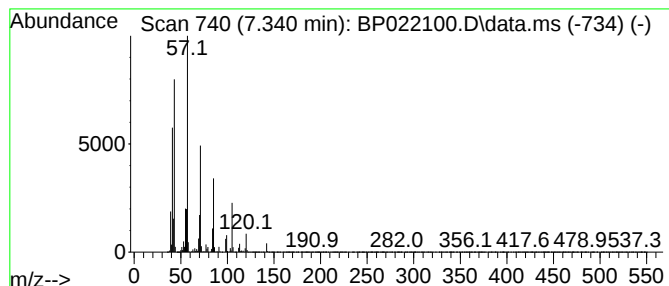
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.340	91.02 ng/ul	5825620	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	93
2			Decane	142	C10H22	000124-18-5	90
3			Decane	142	C10H22	000124-18-5	81
4			Hexadecane	226	C16H34	000544-76-3	53
5			Tridecane	184	C13H28	000629-50-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
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Misc :
ALS Vial : 6 Sample Multiplier: 1

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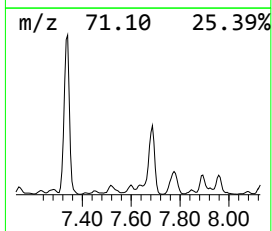
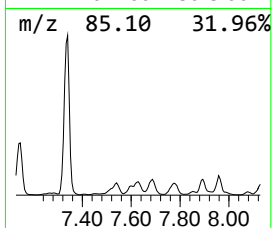
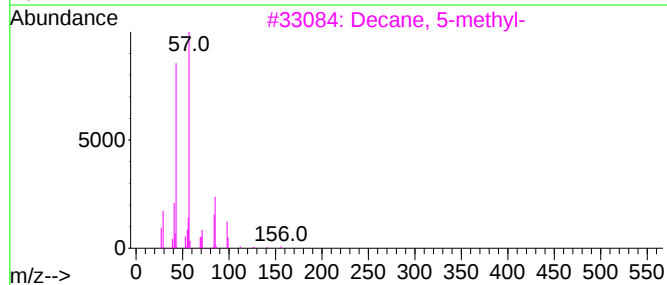
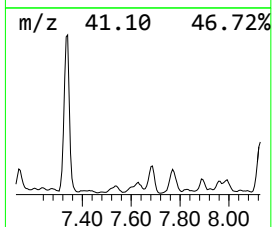
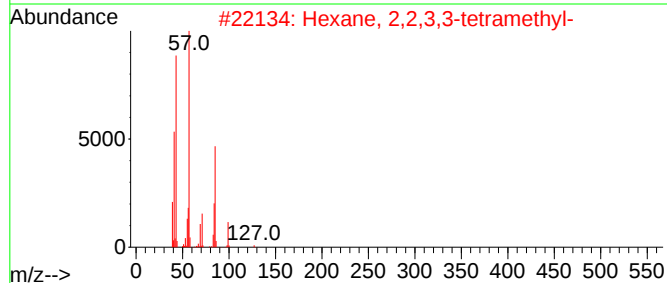
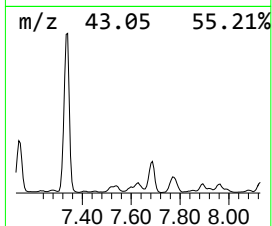
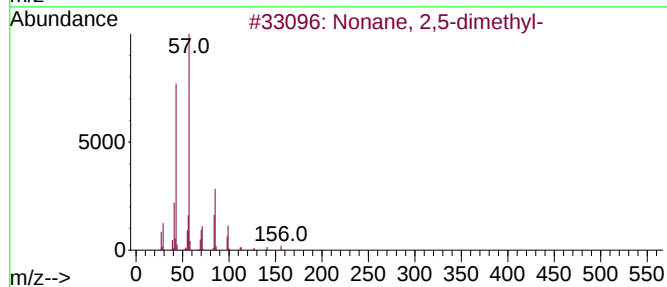
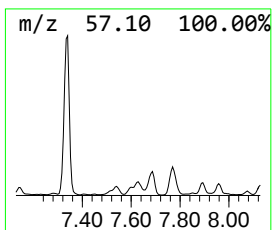
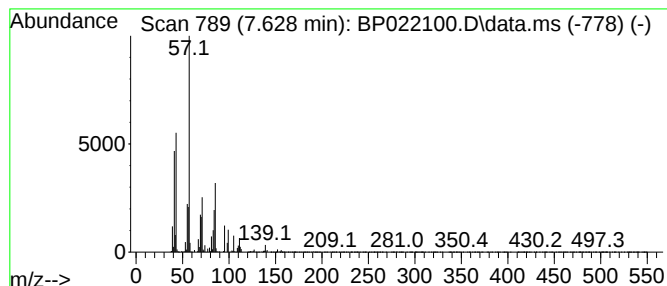
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.628	8.77 ng/ul	561012	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 2,5-dimethyl-	156	C11H24	017302-27-1	53
2			Hexane, 2,2,3,3-tetramethyl-	142	C10H22	013475-81-5	53
3			Decane, 5-methyl-	156	C11H24	013151-35-4	52
4			Pentane, 3-ethyl-2,3-dimethyl-	128	C9H20	016747-33-4	47
5			Carbonic acid, decyl prop-1-en-2...	242	C14H26O3	1000382-90-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

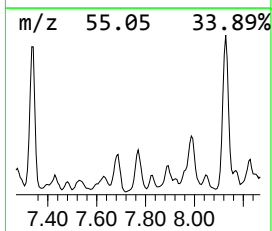
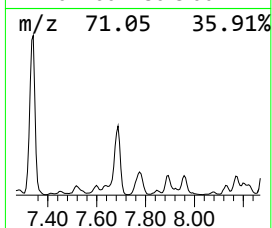
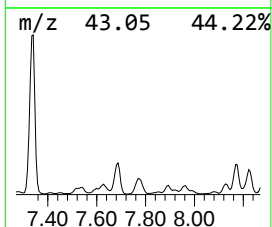
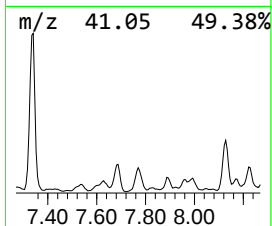
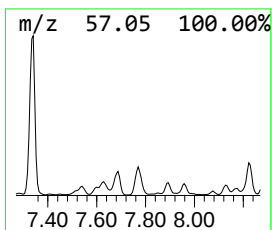
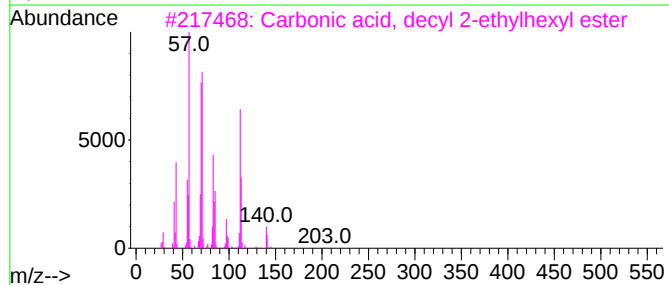
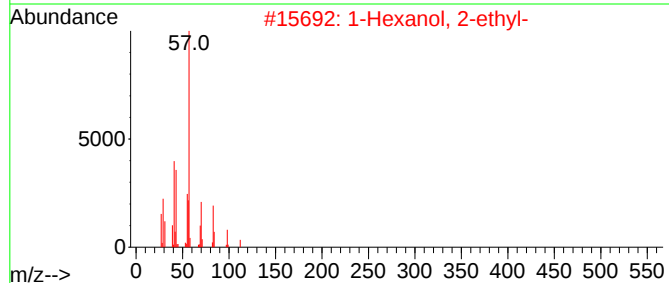
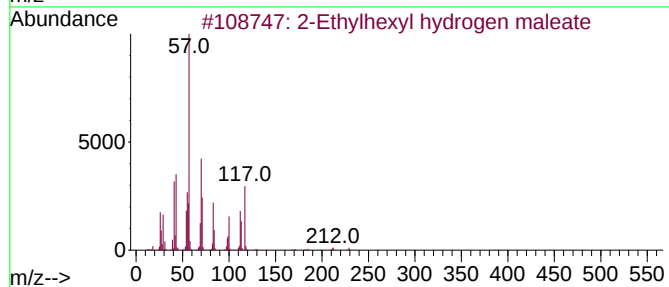
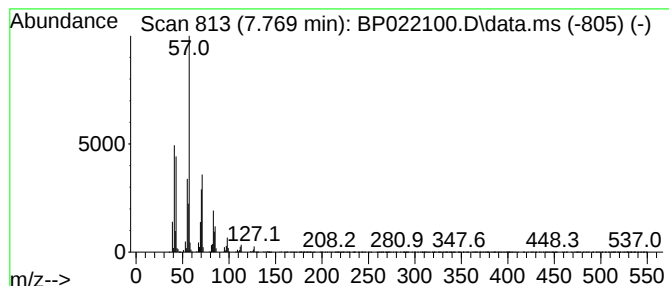
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 2-Ethylhexyl hydrogen maleate Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD			R.T.
7.769	11.99 ng/ul	767510	1,4-Dichlorobenzene-d4			7.699
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Ethylhexyl hydrogen maleate		228	C12H20O4	002370-71-0	59
2	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	59
3	Carbonic acid, decyl 2-ethylhexy...		314	C19H38O3	1000383-13-7	59
4	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	53
5	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

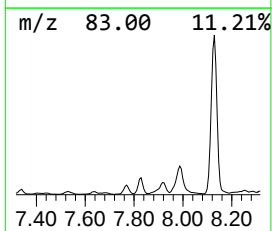
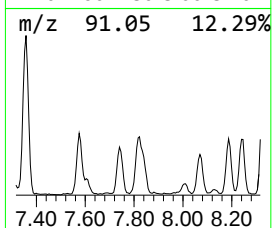
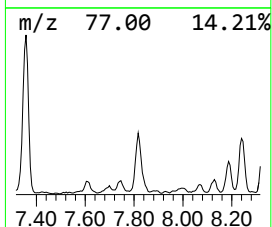
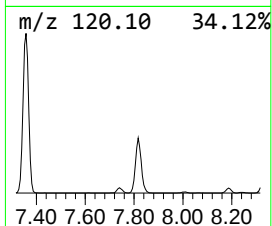
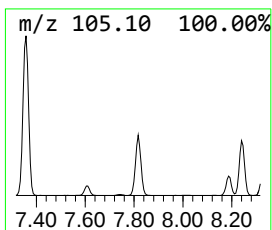
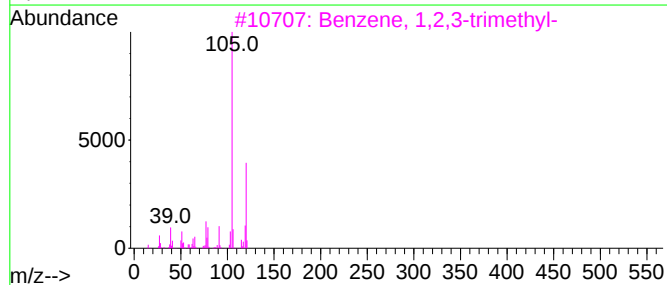
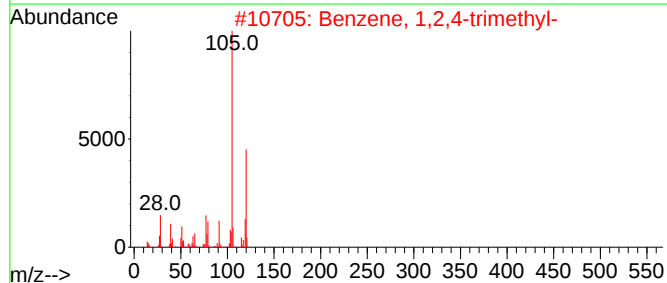
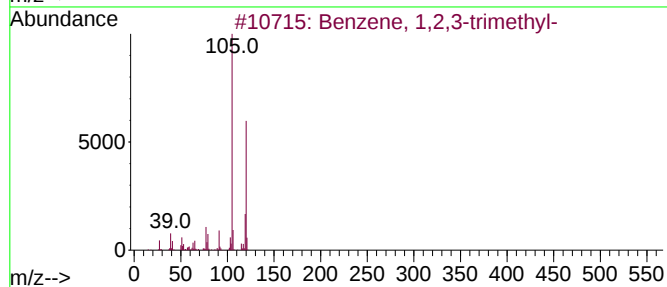
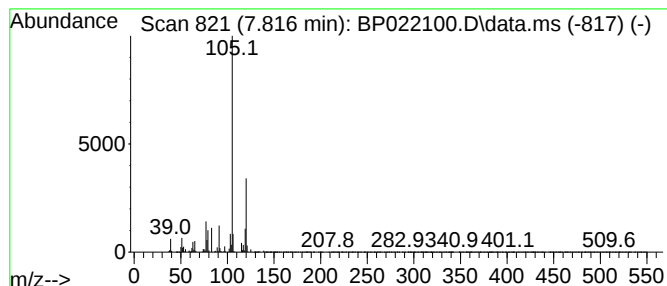
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzene, 1,2,4-trimethyl- Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD			R.T.
7.816	13.10 ng/ul	838166	1,4-Dichlorobenzene-d4			7.699
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-		120	C9H12	000526-73-8	91
2	Benzene, 1,2,4-trimethyl-		120	C9H12	000095-63-6	91
3	Benzene, 1,2,3-trimethyl-		120	C9H12	000526-73-8	87
4	Mesitylene		120	C9H12	000108-67-8	87
5	Mesitylene		120	C9H12	000108-67-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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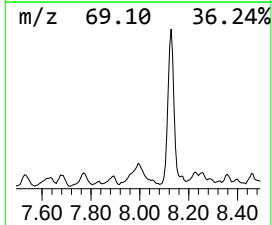
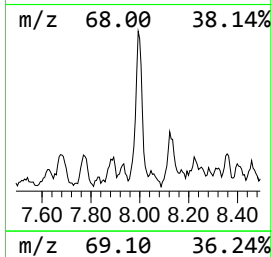
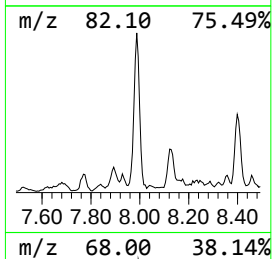
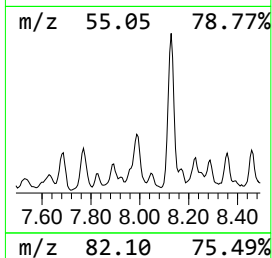
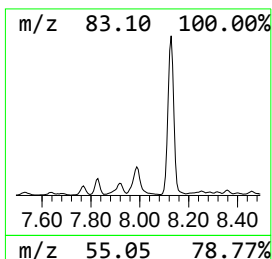
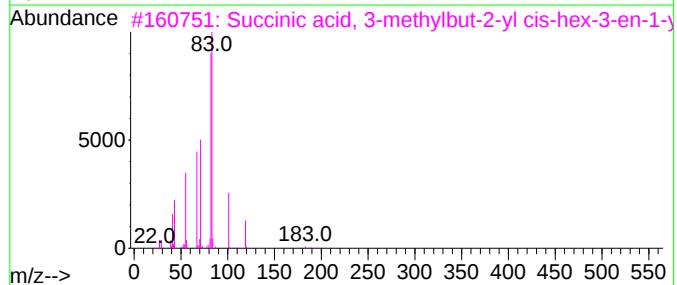
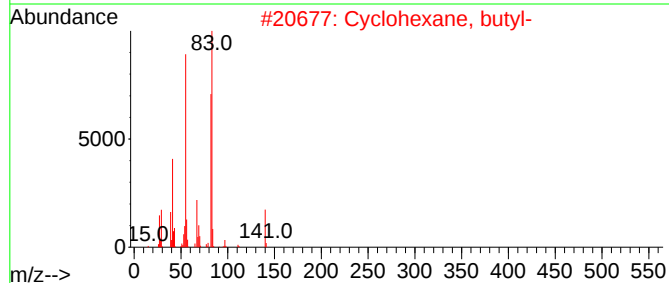
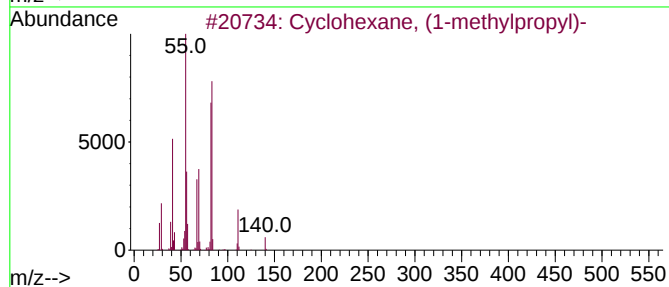
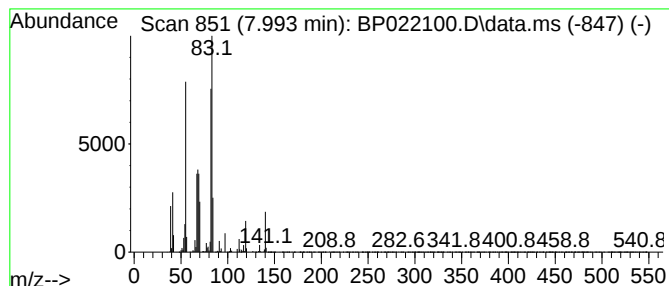
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Cyclic7.993 Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.993	8.32 ng/ul	532355	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (1-methylpropyl)-	140	C10H20	007058-01-7	53
2			Cyclohexane, butyl-	140	C10H20	001678-93-9	53
3			Succinic acid, 3-methylbut-2-yl ...	270	C15H26O4	1000391-08-2	50
4			Succinic acid, 3-methylbut-2-yl ...	270	C15H26O4	1000391-11-1	50
5			Cyclopentane, 1-methyl-3-(1-meth...	126	C9H18	053771-88-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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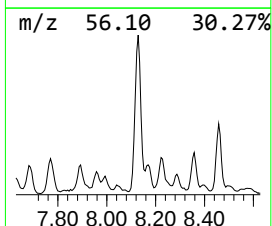
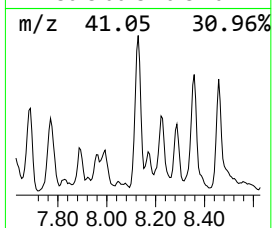
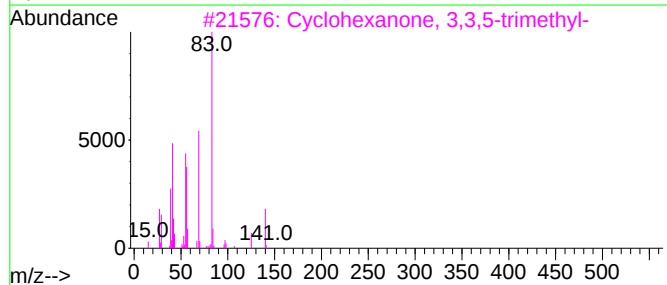
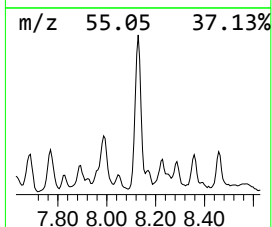
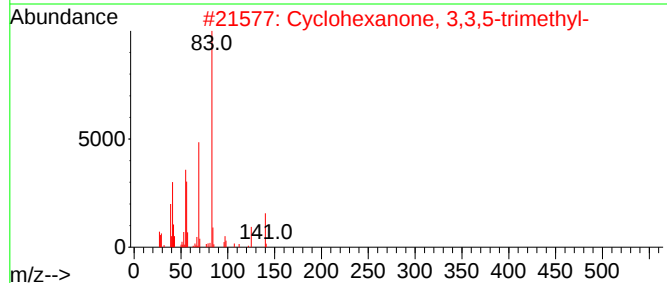
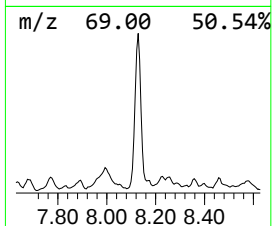
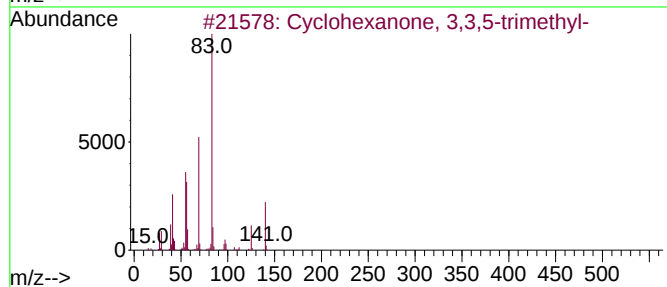
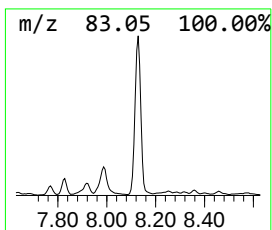
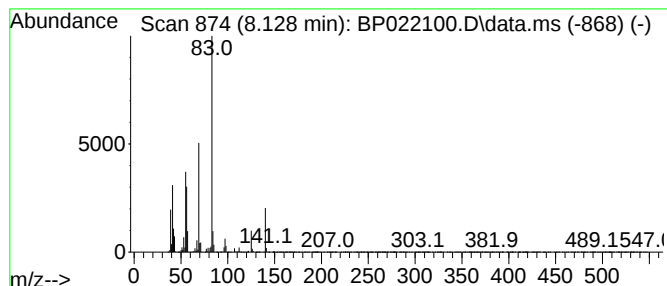
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Cyclohexanone, 3,3,5-trimet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.128	33.28 ng/ul	2129730	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	97
2			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	95
3			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	94
4			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
5			Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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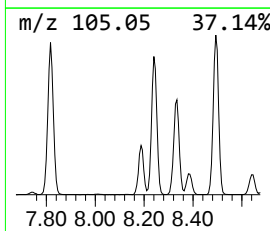
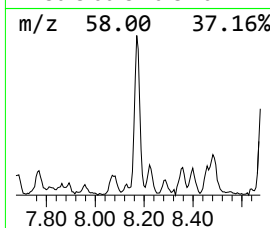
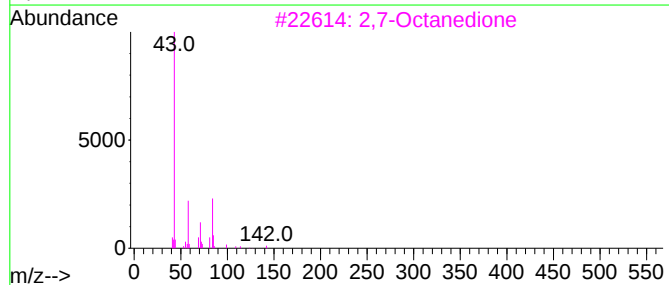
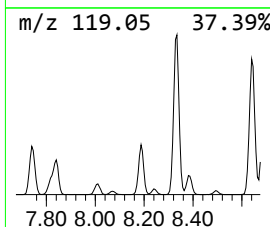
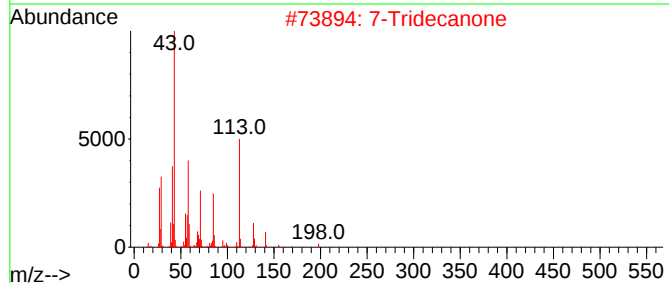
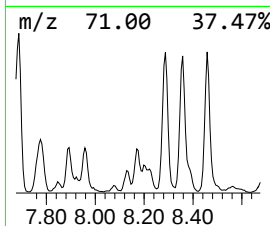
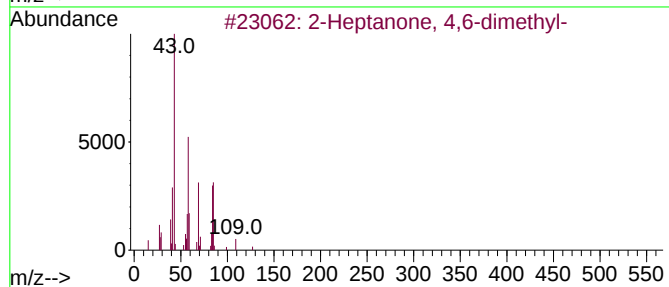
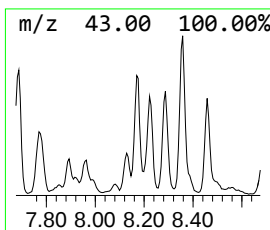
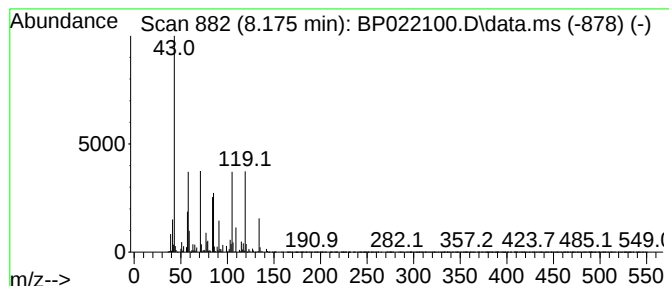
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 unknown-01 Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.175	7.29 ng/ul	466535	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 4,6-dimethyl-			142	C9H18O	019549-80-5	38
2	7-Tridecanone			198	C13H26O	000462-18-0	25
3	2,7-Octanedione			142	C8H14O2	001626-09-1	22
4	7-Tridecanone			198	C13H26O	000462-18-0	17
5	Hexane, 3-ethyl-2-methyl-			128	C9H20	016789-46-1	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
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Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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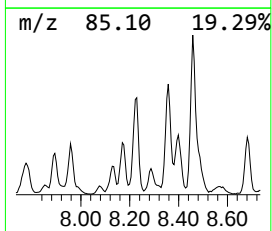
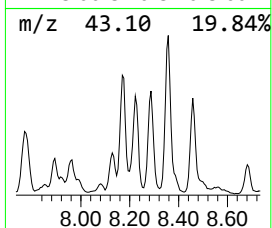
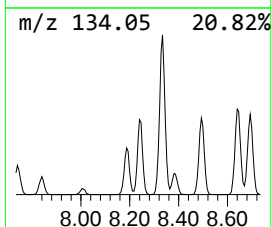
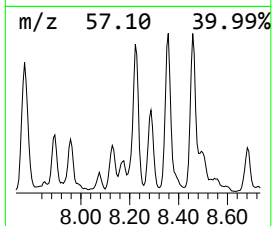
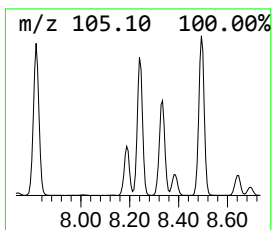
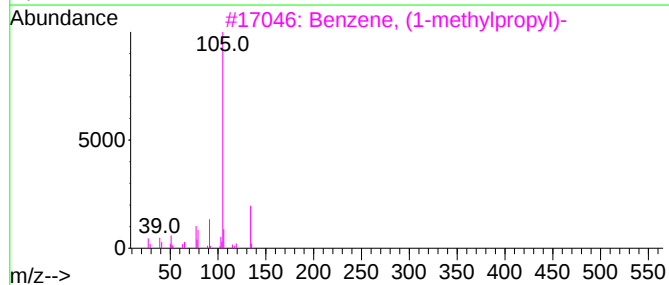
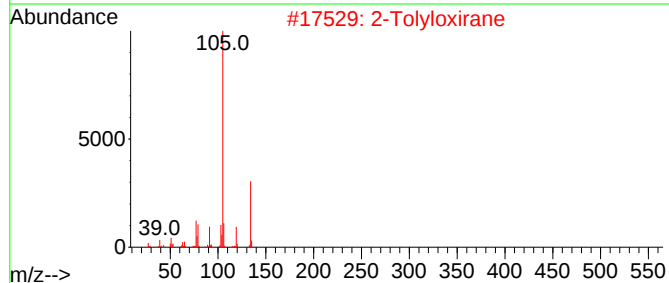
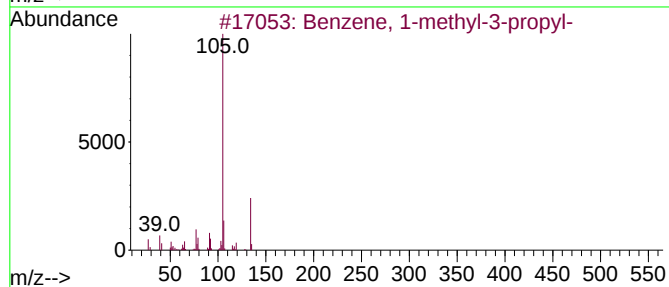
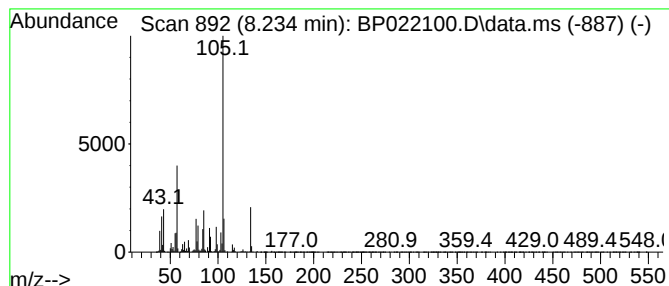
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Benzene, 1-methyl-3-propyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.234	16.10 ng/ul	1030400	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	60
2			2-Tolyloxirane	134	C9H10O	002783-26-8	55
3			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	55
4			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	55
5			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
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ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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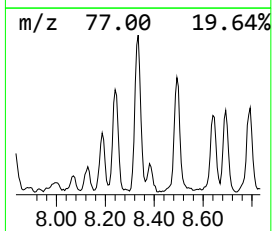
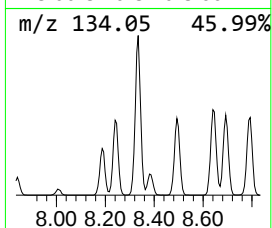
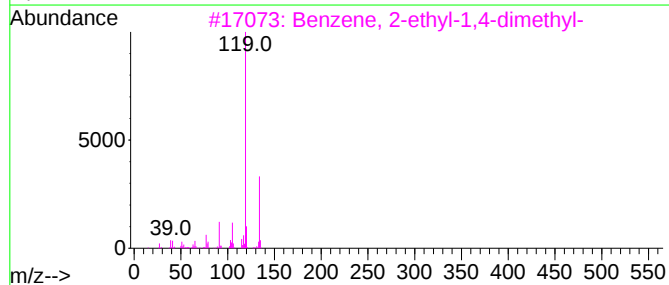
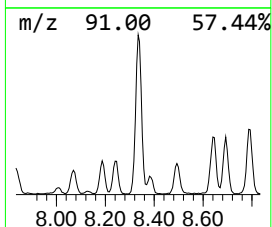
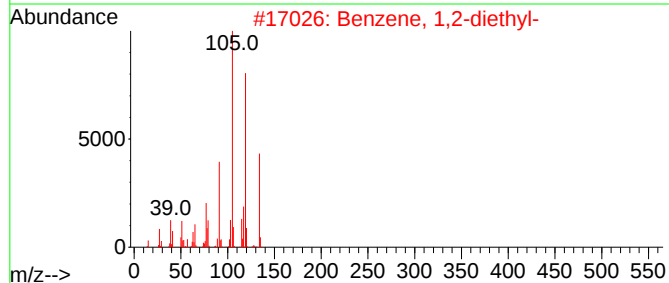
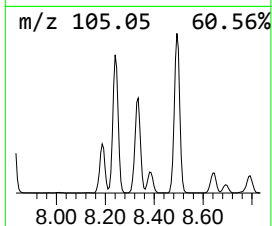
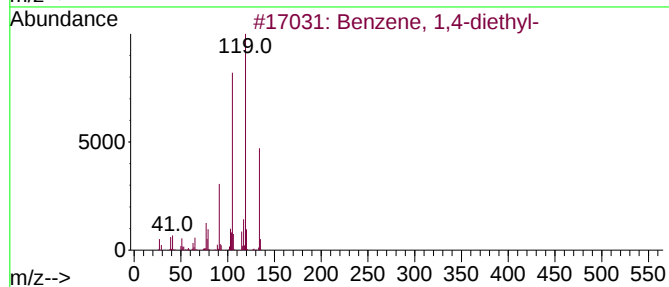
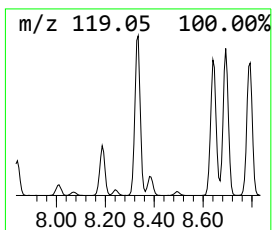
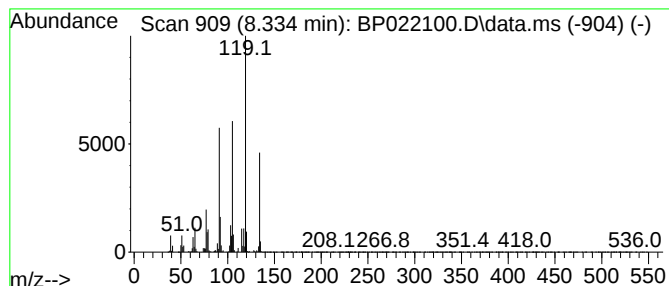
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Benzene, 1,4-diethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.334	19.03 ng/ul	1218220	1,4-Dichlorobenzene-d4	7.699

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, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

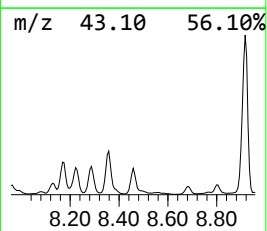
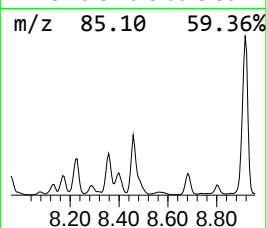
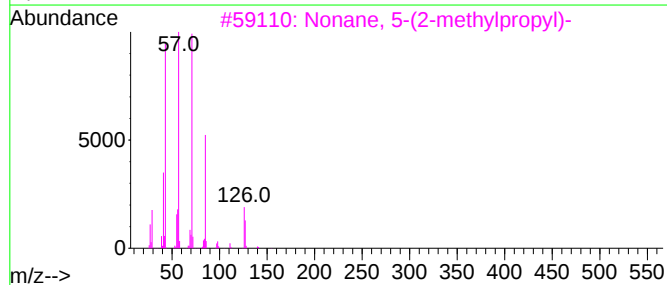
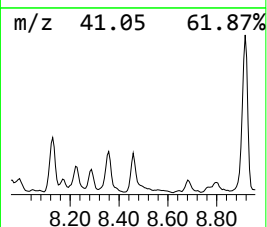
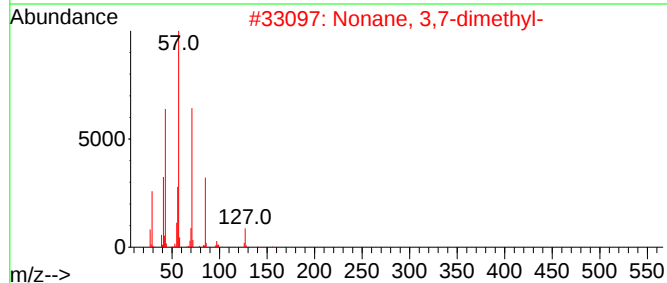
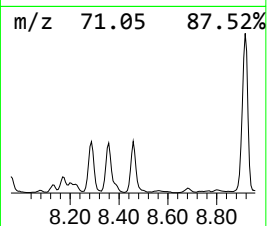
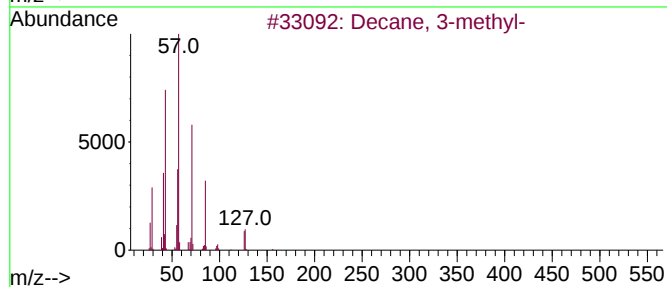
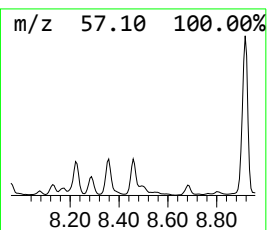
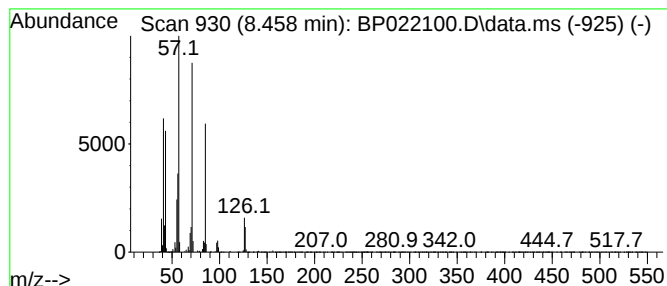
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.458	14.30 ng/ul	915367	1,4-Dichlorobenzene-d4	7.699	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 3-methyl-	156	C11H24	013151-34-3	91
2	Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	87
3	Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	78
4	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	72
5	Nonane, 1-iodo-	254	C9H19I	004282-42-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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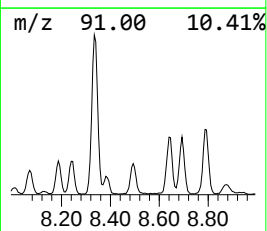
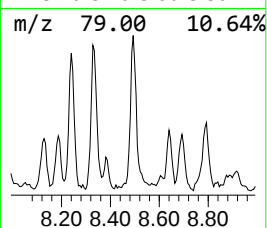
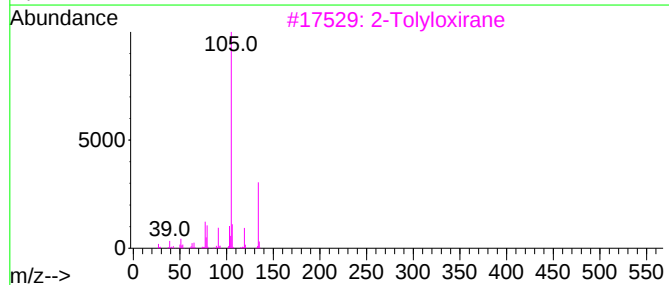
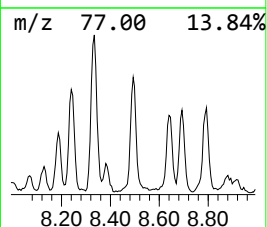
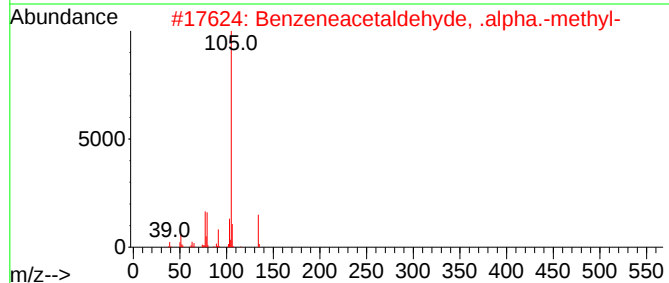
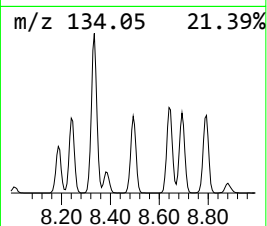
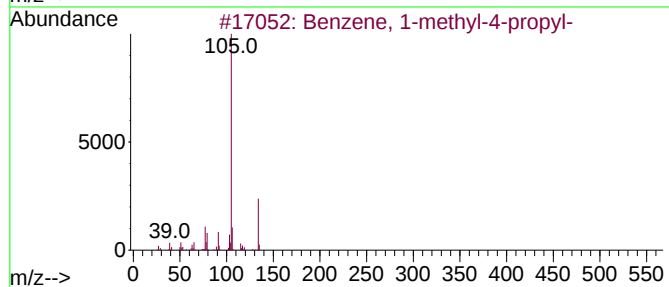
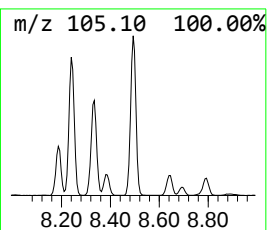
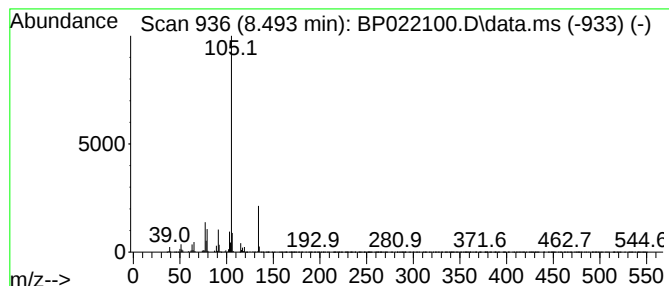
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Benzene, 1-methyl-4-propyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.493	11.90 ng/ul	761872	1,4-Dichlorobenzene-d4	7.699

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			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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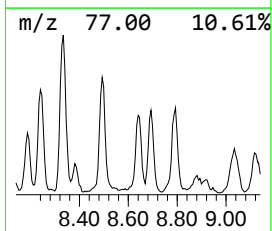
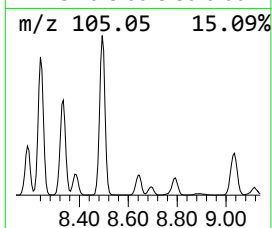
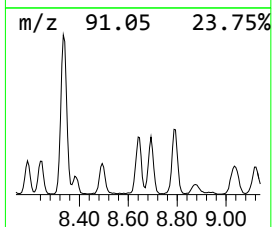
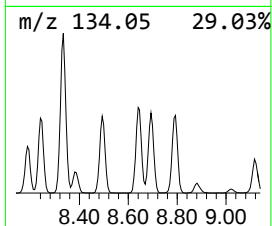
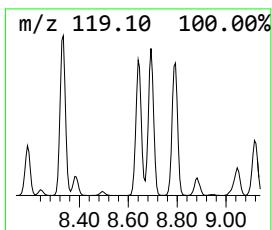
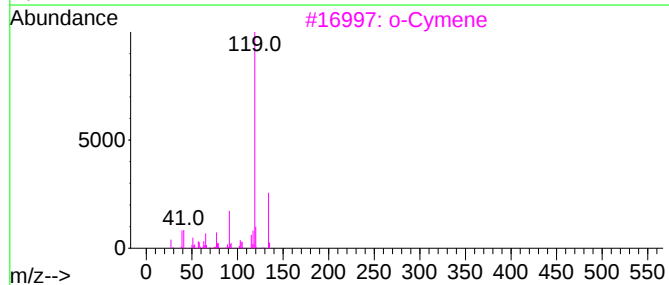
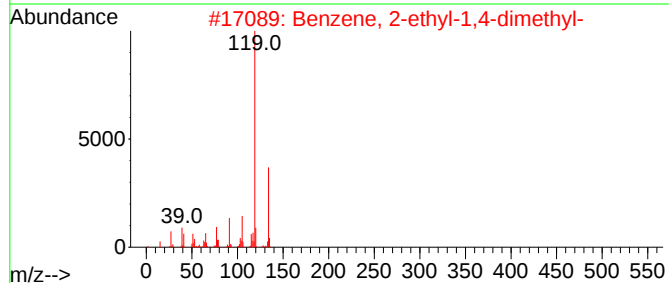
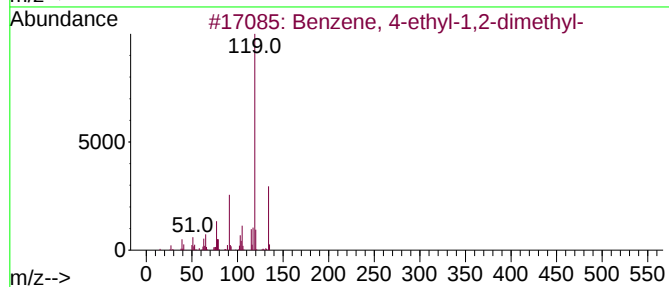
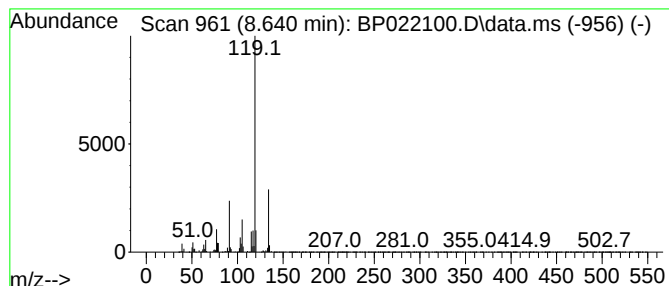
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.640	9.70 ng/ul	620905	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3			o-Cymene	134	C10H14	000527-84-4	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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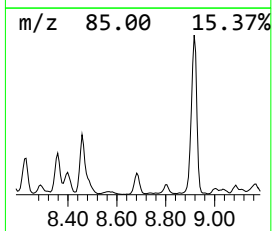
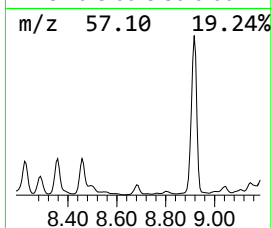
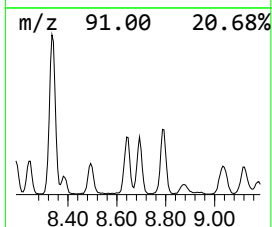
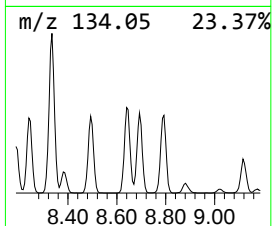
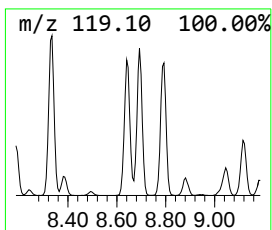
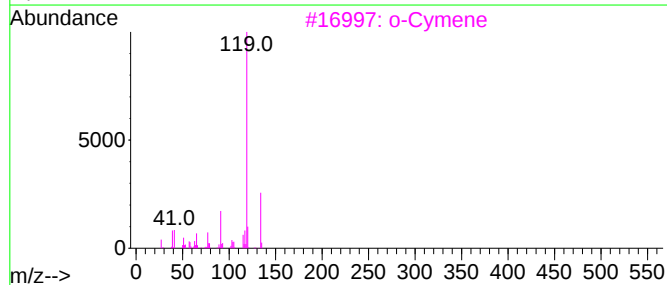
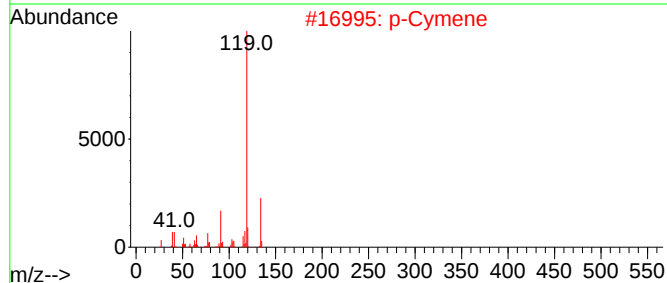
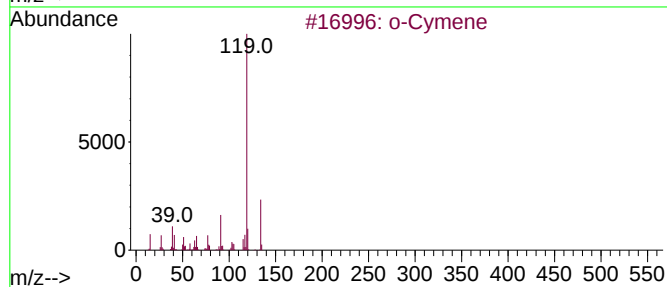
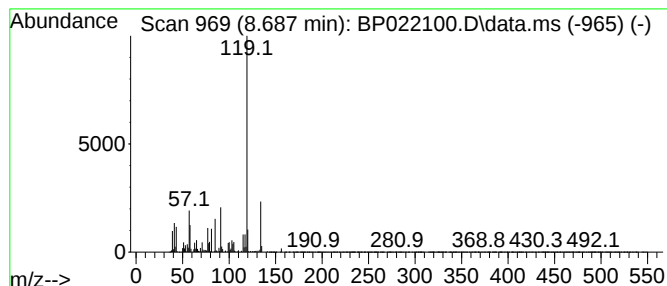
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 o-Cymene Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.687	15.54 ng/ul	994802	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	95
2			p-Cymene	134	C10H14	000099-87-6	95
3			o-Cymene	134	C10H14	000527-84-4	95
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5			o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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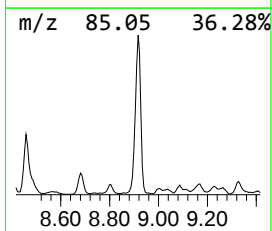
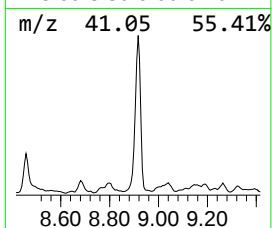
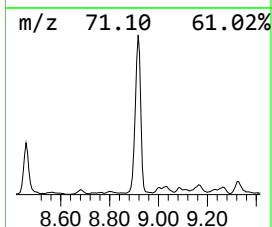
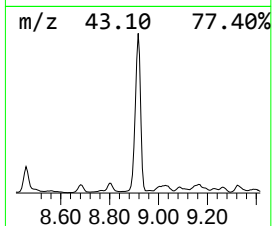
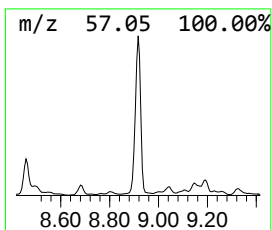
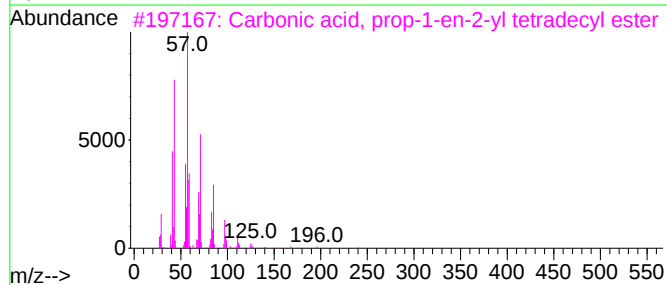
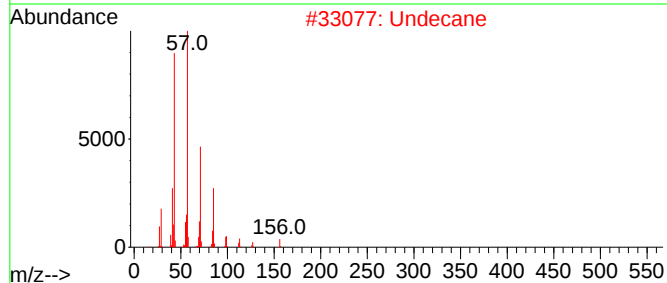
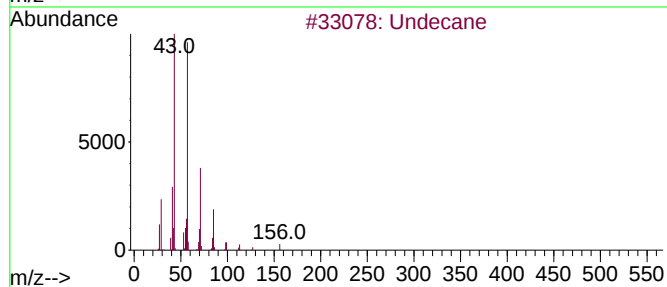
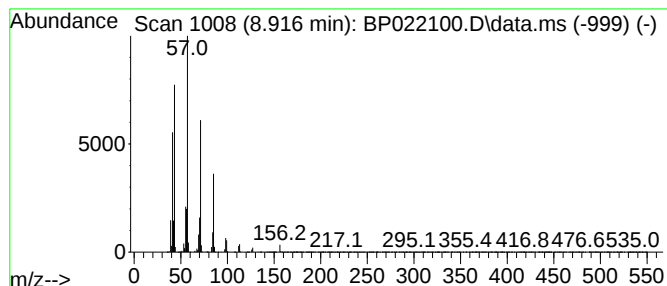
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.916	63.43 ng/ul	4059680	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	87
2	Undecane			156	C11H24	001120-21-4	86
3	Carbonic acid, prop-1-en-2-yl te...			298	C18H34O3	1000382-90-9	83
4	Hexadecane			226	C16H34	000544-76-3	80
5	Tetradecane			198	C14H30	000629-59-4	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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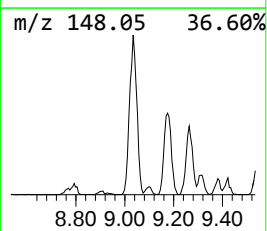
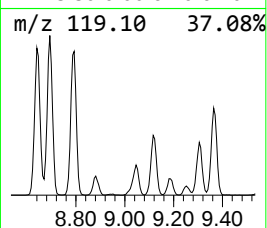
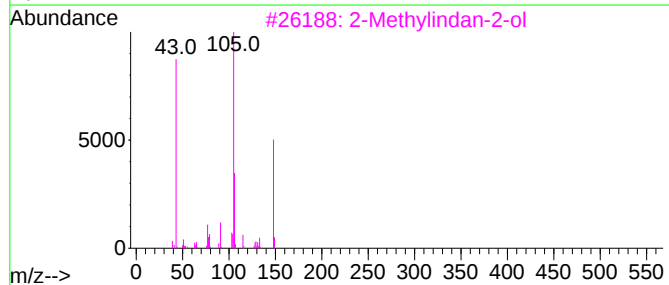
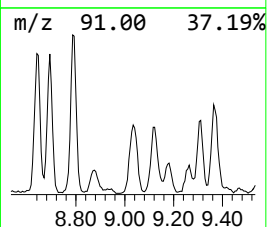
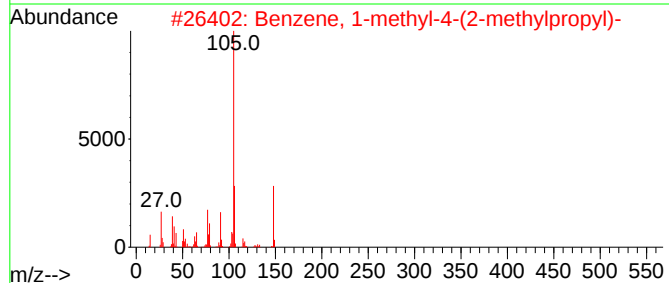
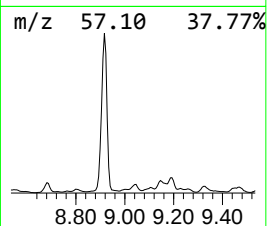
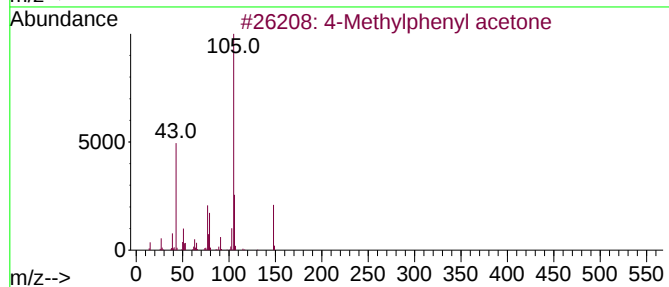
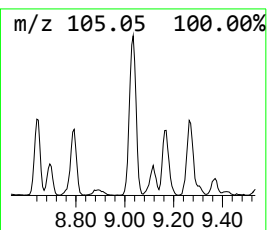
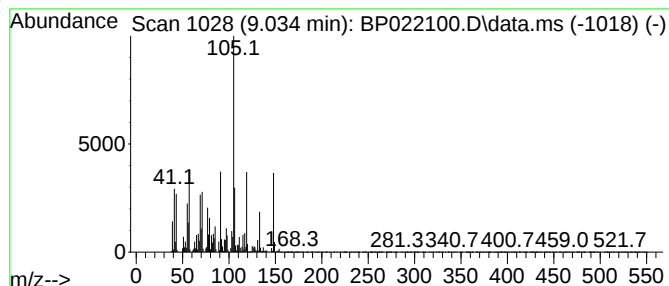
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 4-Methylphenyl acetone Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.034	14.68 ng/ul	939770	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Methylphenyl acetone	148	C10H12O	002096-86-8	55
2			Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	49
3			2-Methylindan-2-ol	148	C10H12O	033223-84-6	43
4			Benzene, (1-methylbutyl)-	148	C11H16	002719-52-0	42
5			3-Buten-2-ol, 4-phenyl-, (E)-	148	C10H12O	1000163-70-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC75ME

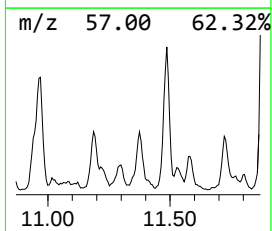
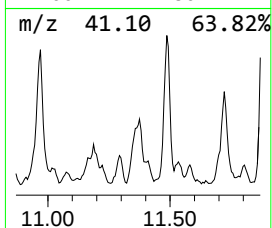
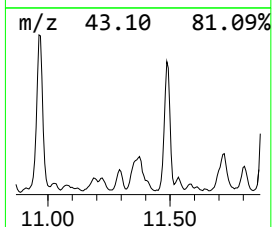
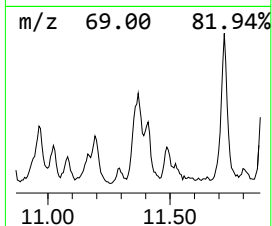
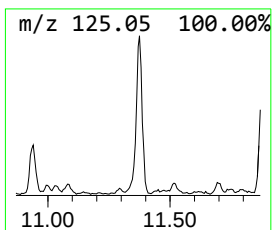
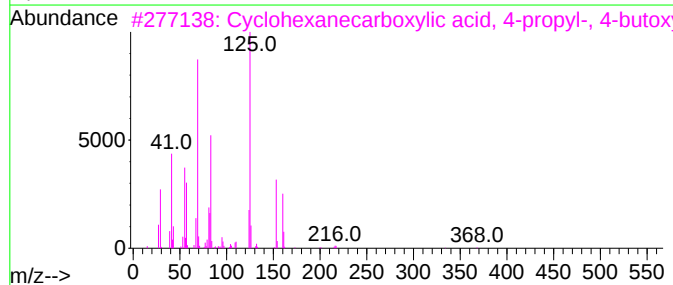
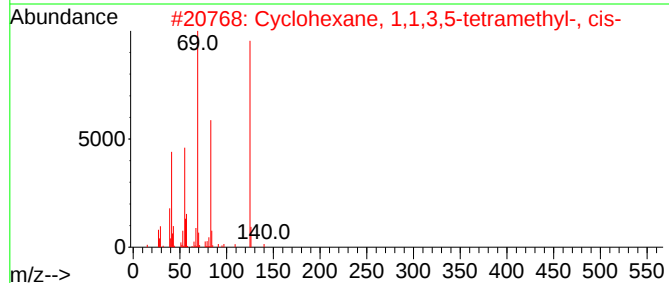
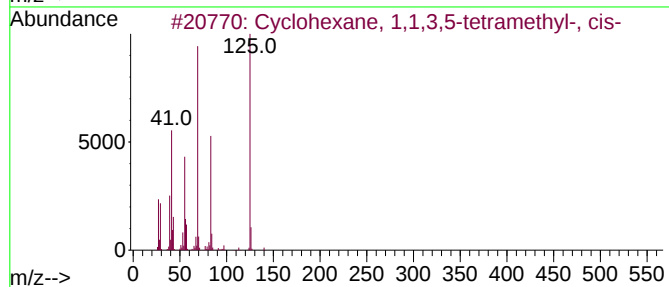
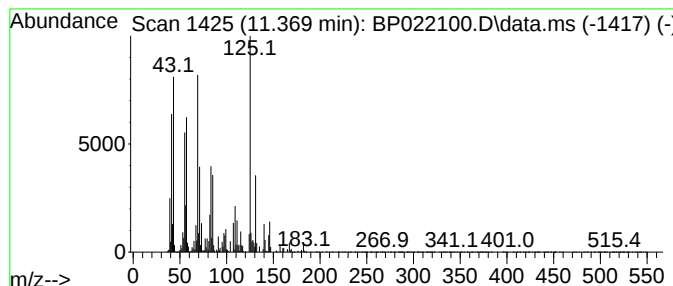
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 (DEL) Alkane: Cyclic11.369 Concentration Rank 62

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.369	2.14 ng/ul	795126	Naphthalene-d8	10.458

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-32-9	47
2			Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-32-9	47
3			Cyclohexanecarboxylic acid, 4-pr...	368	C22H28N2O3	075941-67-2	47
4			Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-31-8	47
5			Cyclohexane, 1,2-diethyl-1-methyl-	154	C11H22	061141-79-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

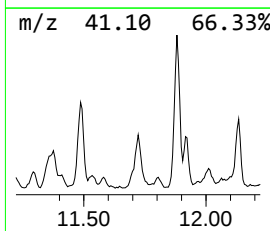
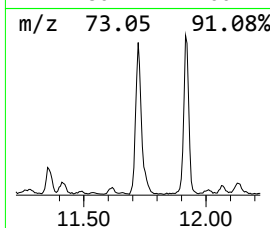
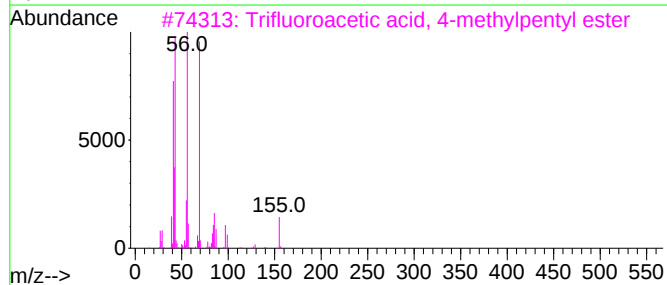
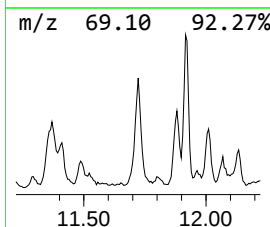
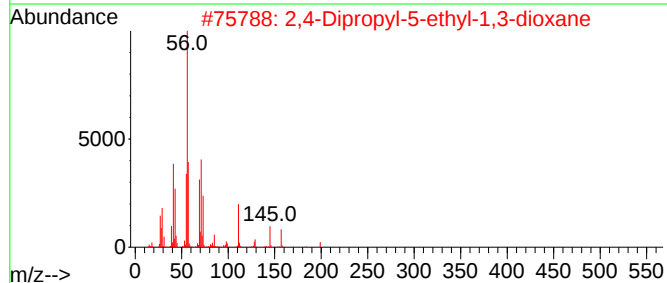
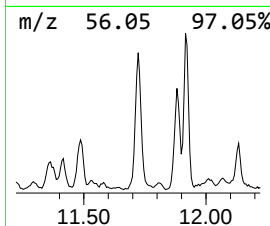
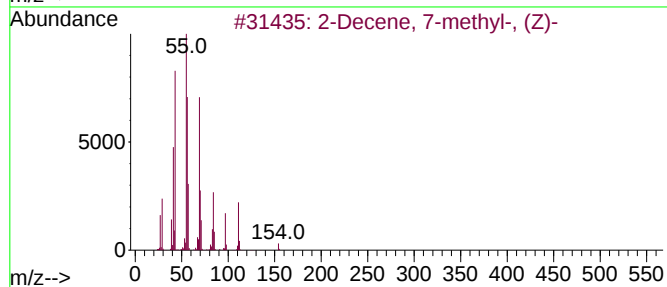
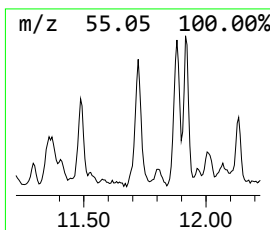
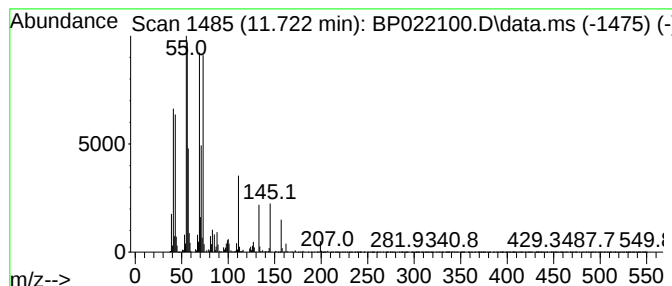
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 61

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.722	2.29 ng/ul	851214	Naphthalene-d8	10.458	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Decene, 7-methyl-, (Z)-	154	C11H22	074630-23-2	47
2	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	38
3	Trifluoroacetic acid, 4-methylpe...	198	C8H13F3O2	000327-71-9	38
4	1-Octene, 2,7-dimethyl-	140	C10H20	033718-03-5	38
5	4-Undecene, 10-methyl-, (E)-	168	C12H24	074630-60-7	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC75ME

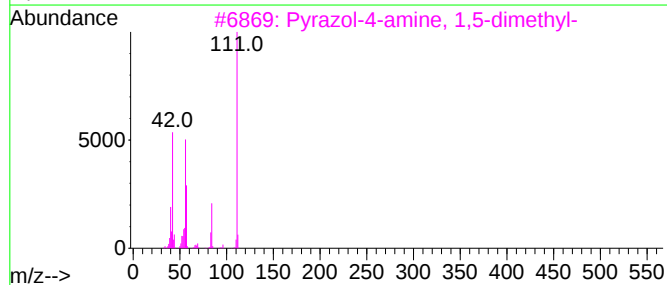
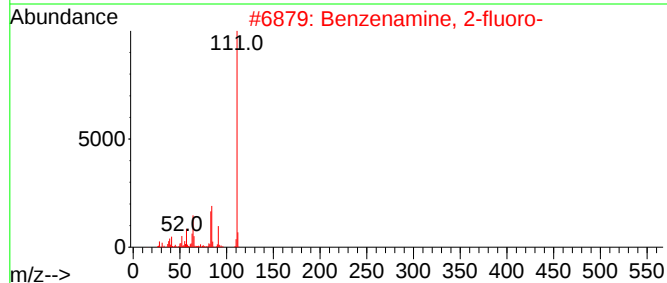
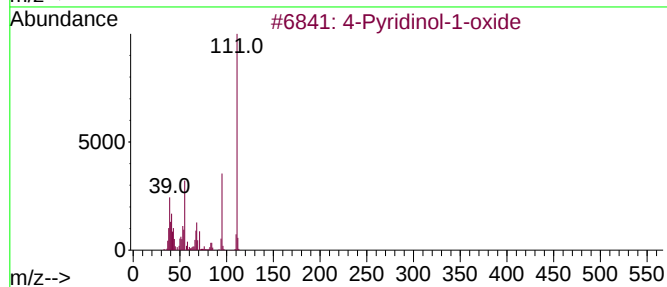
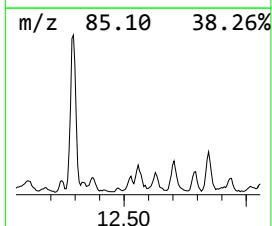
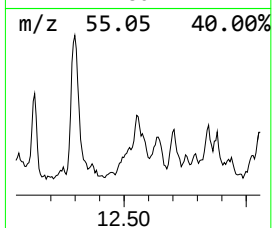
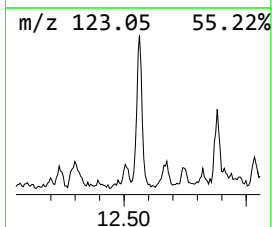
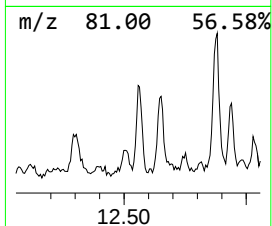
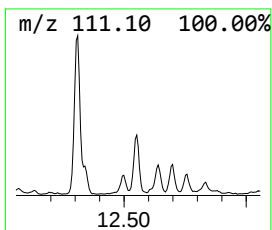
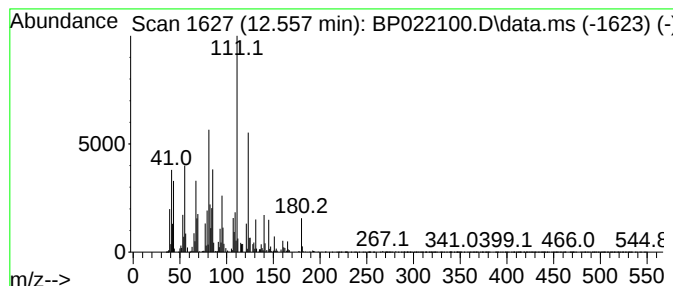
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 unknown-02 Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.557	7.40 ng/ul	561272	Acenaphthene-d10	14.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Pyridinol-1-oxide	111	C5H5NO2	006890-62-6	30
2			Benzenamine, 2-fluoro-	111	C6H6FN	000348-54-9	27
3			Pyrazol-4-amine, 1,5-dimethyl-	111	C5H9N3	121983-36-6	27
4			Cytosine	111	C4H5N3O	000071-30-7	25
5			2-Amino-4-acetamino anisole	180	C9H12N2O2	006375-47-9	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

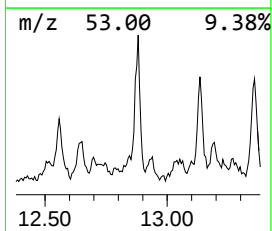
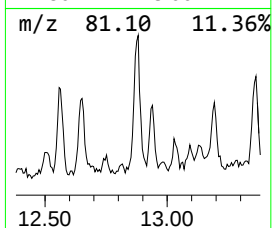
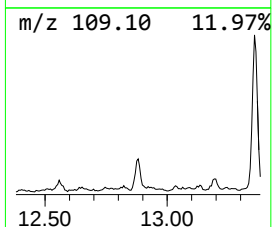
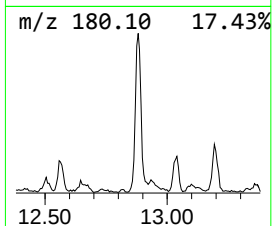
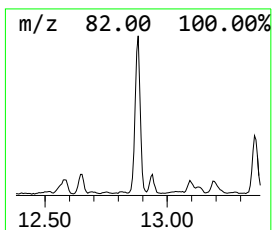
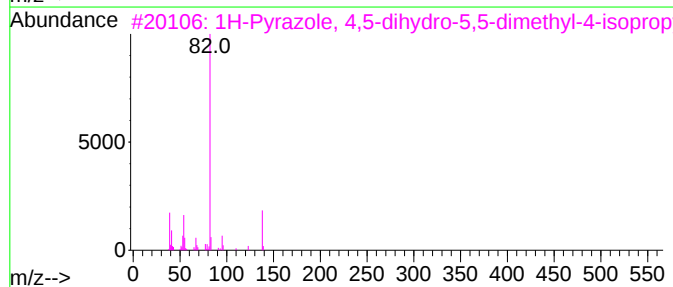
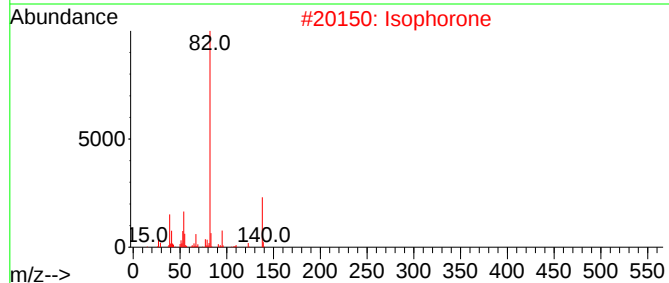
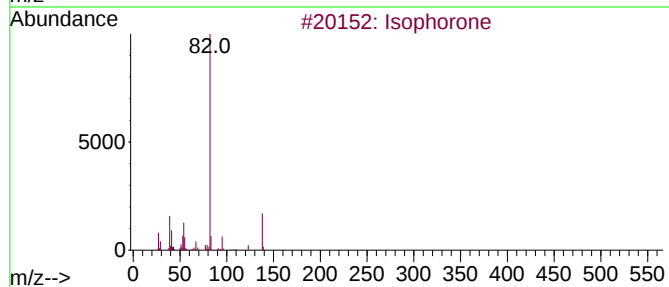
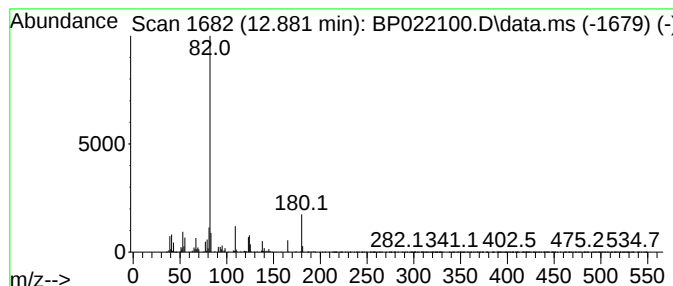
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 32 unknown-03 Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.881	6.57 ng/ul	498566	Acenaphthene-d10		14.310	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Isophorone		138	C9H14O	000078-59-1	50
2	Isophorone		138	C9H14O	000078-59-1	50
3	1H-Pyrazole, 4,5-dihydro-5,5-dim...		138	C8H14N2	106251-09-6	47
4	Isophorone		138	C9H14O	000078-59-1	47
5	1H-Pyrazole, 3-methyl-		82	C4H6N2	001453-58-3	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

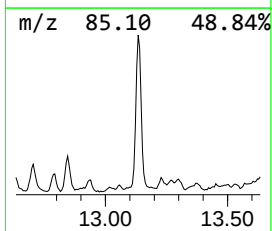
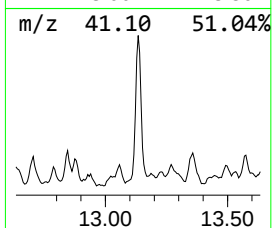
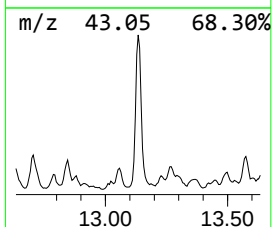
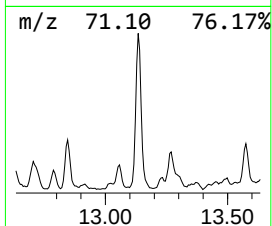
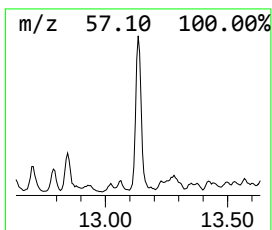
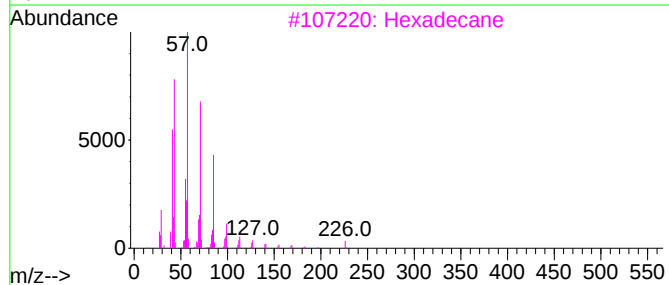
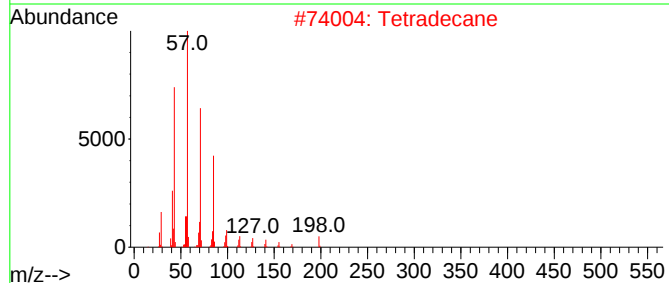
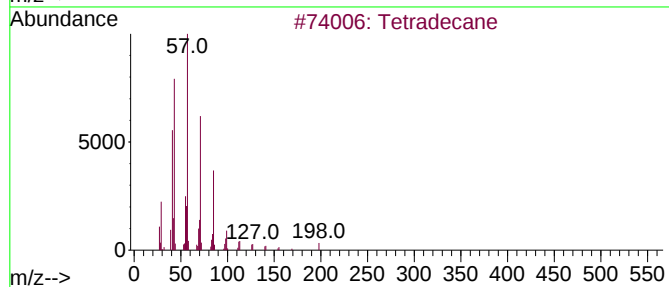
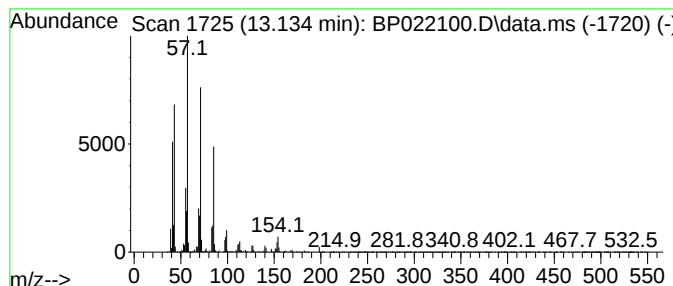
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.134	21.38 ng/ul	1621930	Acenaphthene-d10	14.310	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	95
2	Tetradecane	198	C14H30	000629-59-4	91
3	Hexadecane	226	C16H34	000544-76-3	87
4	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	87
5	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

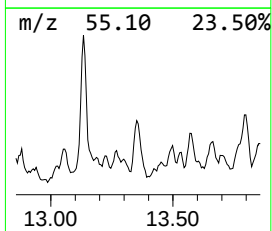
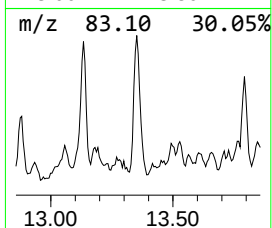
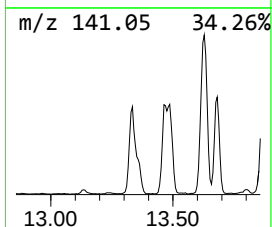
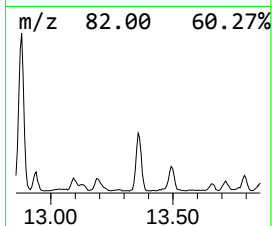
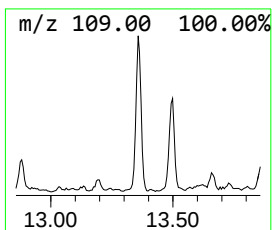
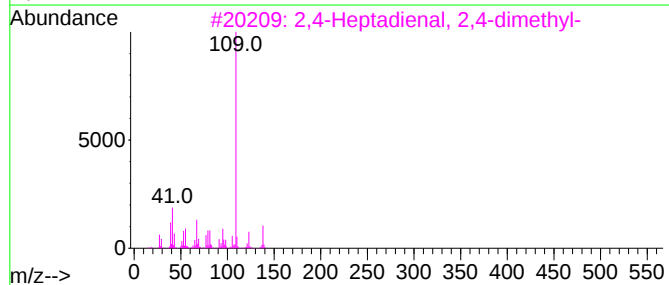
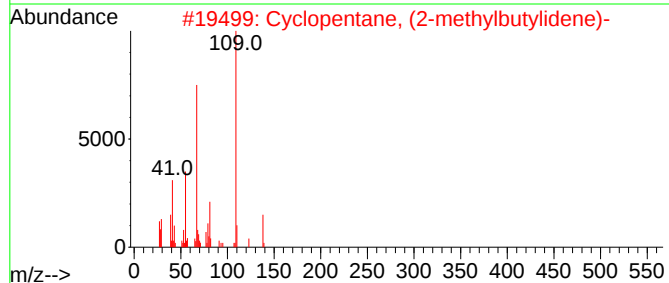
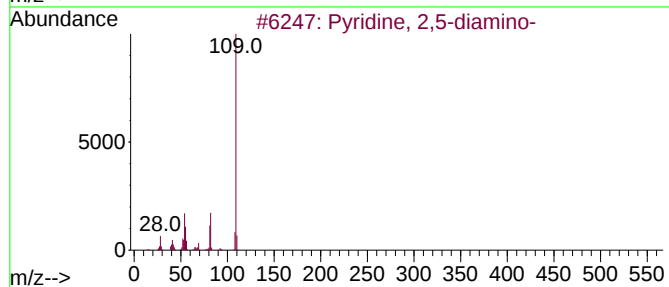
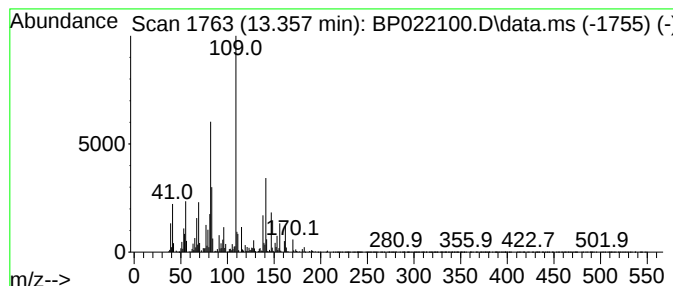
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 35 unknown-04 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.357	15.47 ng/ul	1173810	Acenaphthene-d10	14.310	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyridine, 2,5-diamino-	109	C5H7N3	004318-76-7	49
2	Cyclopentane, (2-methylbutylidene)-	138	C10H18	053366-54-4	43
3	2,4-Heptadienal, 2,4-dimethyl-	138	C9H14O	042452-48-2	43
4	4-Fluorobenzyl mercaptan	142	C7H7FS	015894-04-9	38
5	2,3-Pyridinediamine	109	C5H7N3	000452-58-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

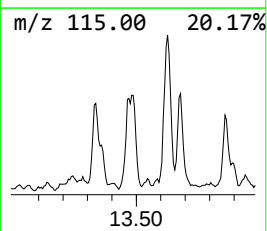
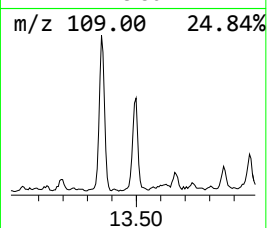
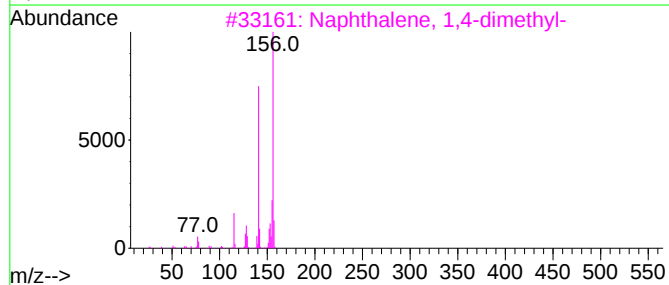
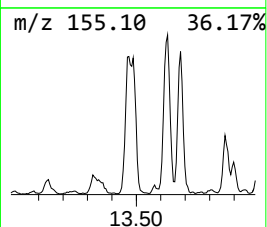
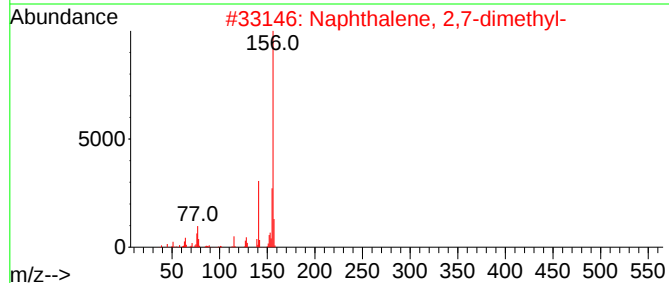
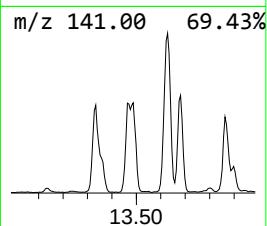
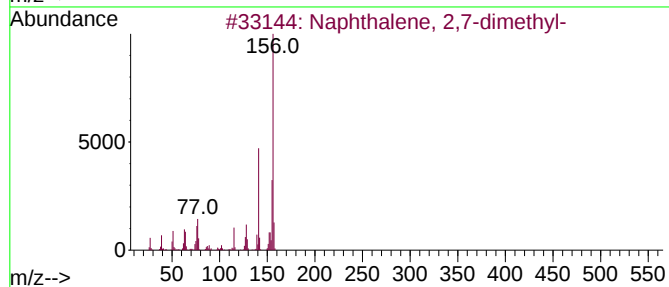
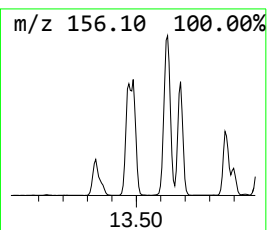
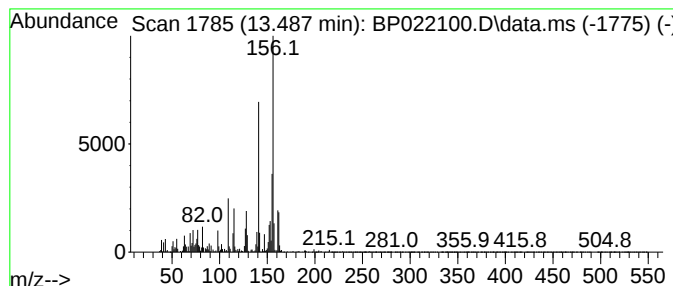
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BNA_P
ClientSampleId :
DDC75ME

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 36 Naphthalene, 2,7-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.487	21.72 ng/ul	1647430	Acenaphthene-d10	14.310	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	94
5	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

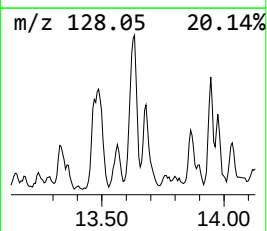
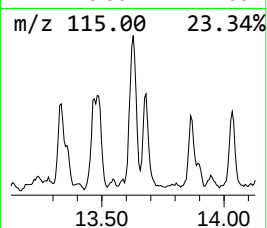
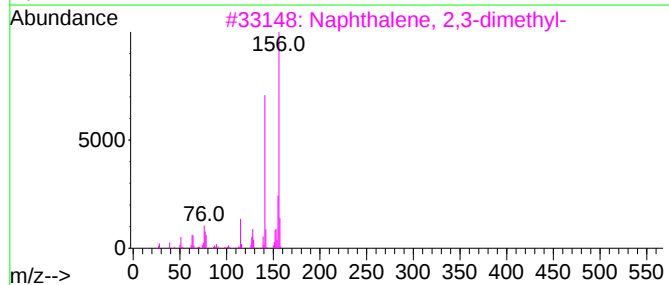
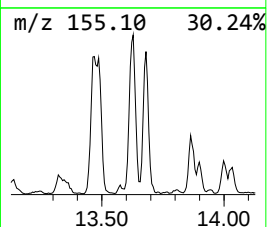
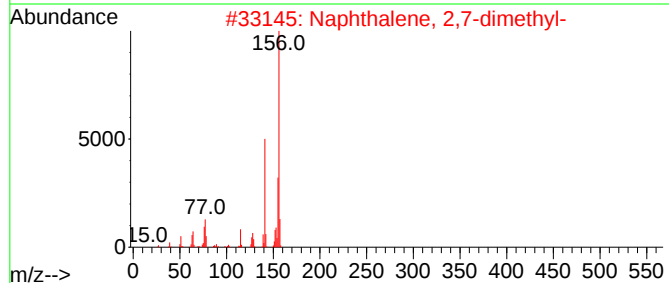
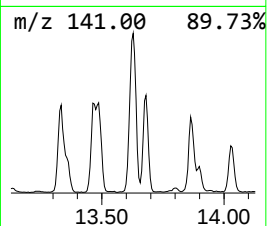
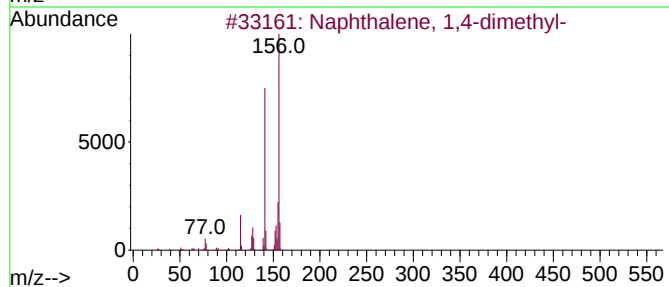
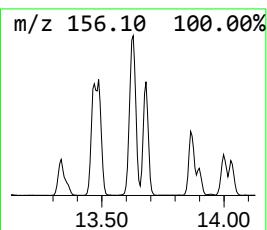
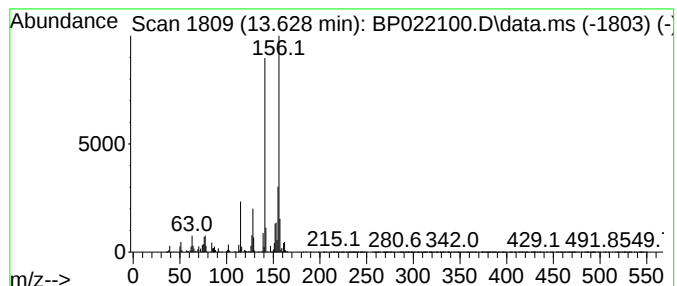
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 37 Naphthalene, 1,4-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.628	16.41 ng/ul	1244530	Acenaphthene-d10	14.310	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
4	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
5	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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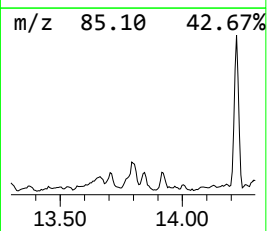
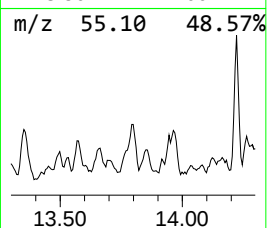
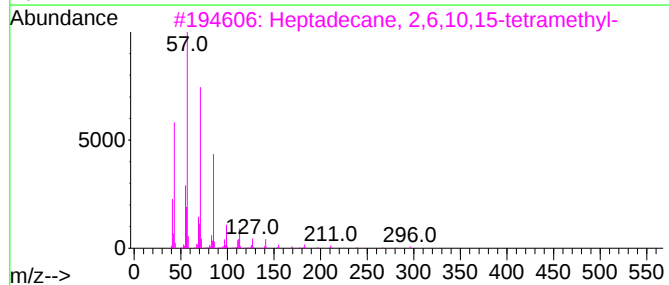
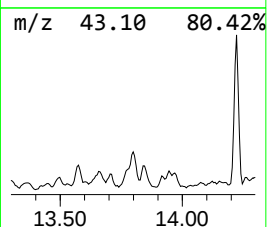
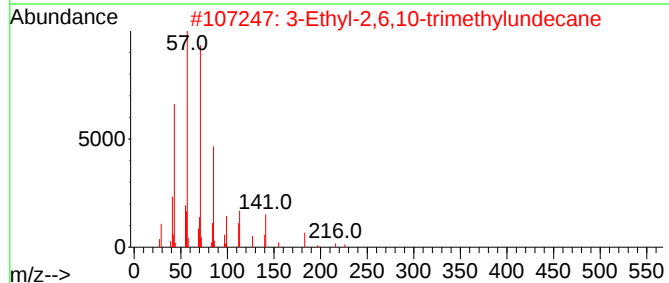
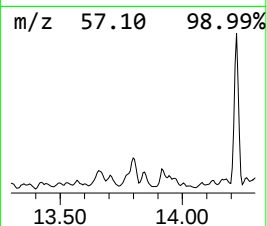
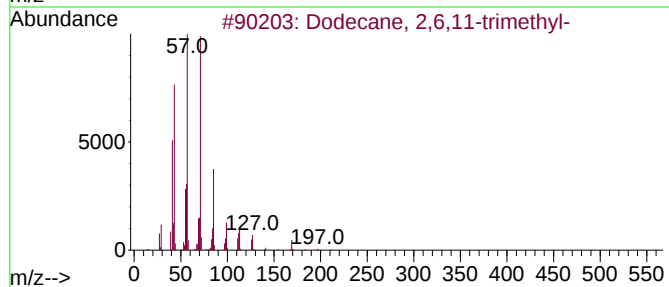
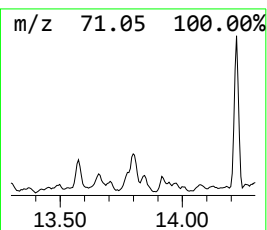
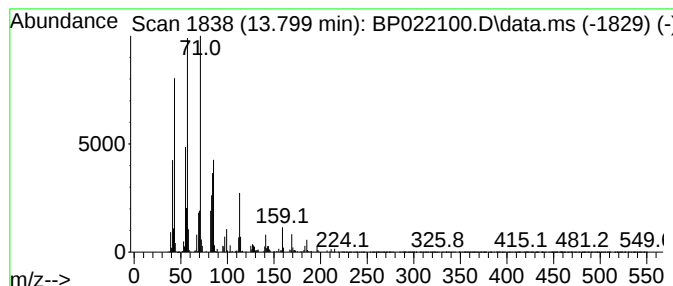
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.799	8.42 ng/ul	638868	Acenaphthene-d10	14.310

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	53
2		3-Ethyl-2,6,10-trimethylundecane	226	C16H34	1000432-25-9	49
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	49
4		Sulfurous acid, decyl pentyl ester	292	C15H32O3S	1000309-14-3	47
5		Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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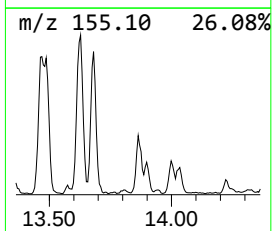
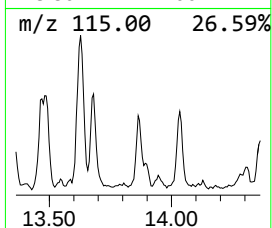
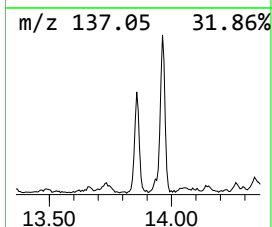
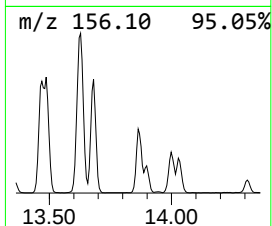
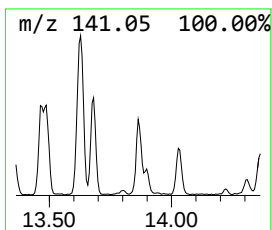
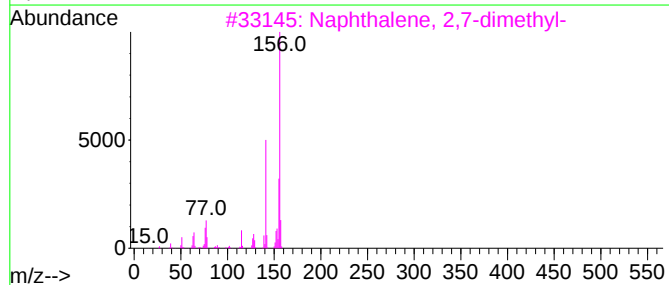
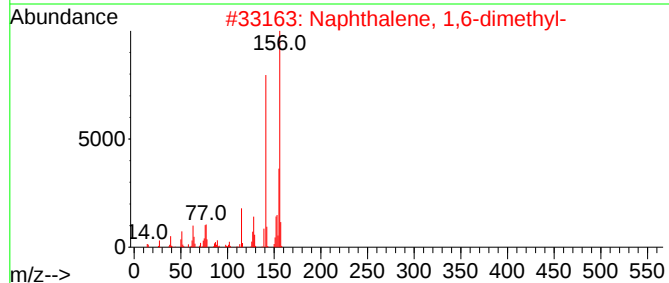
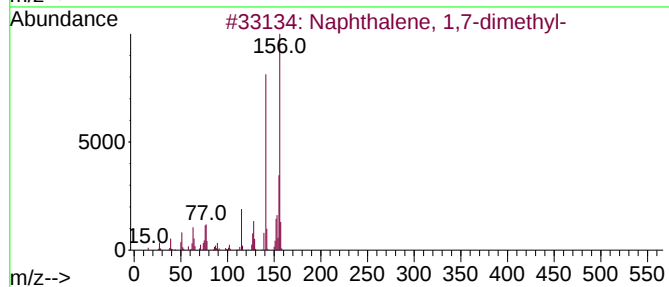
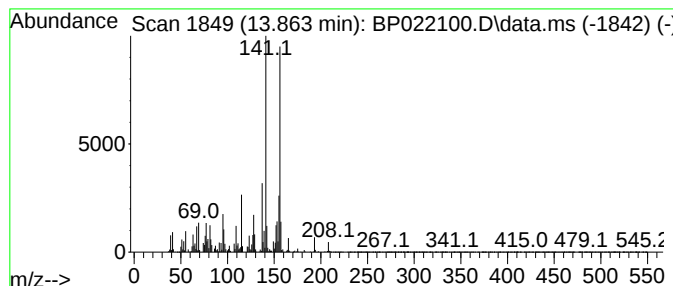
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 Naphthalene, 1,7-dimethyl- Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.863	9.16 ng/ul	694931	Acenaphthene-d10	14.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
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Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

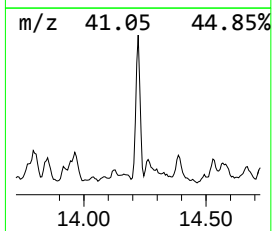
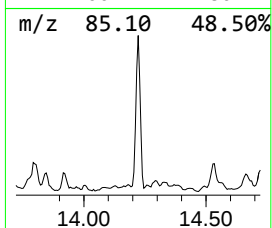
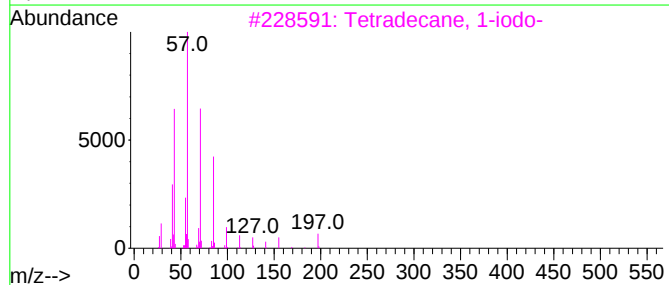
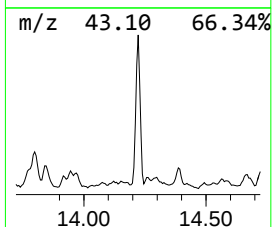
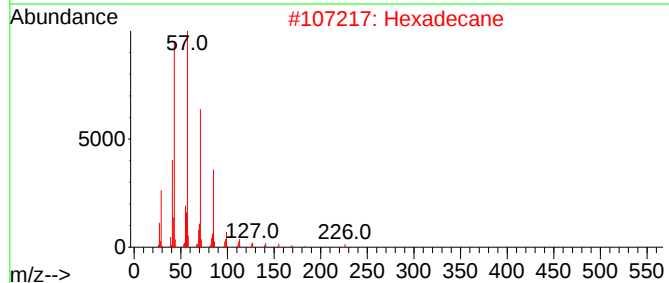
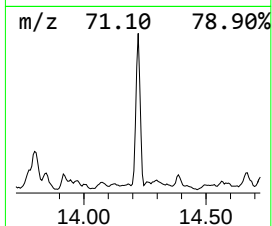
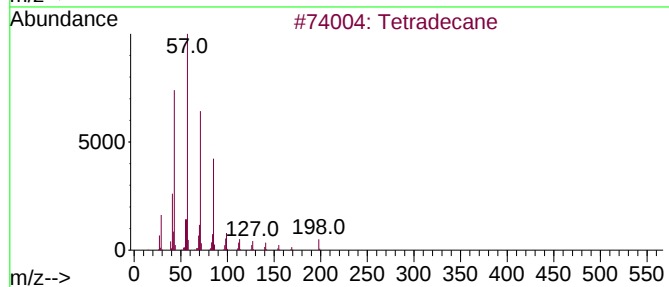
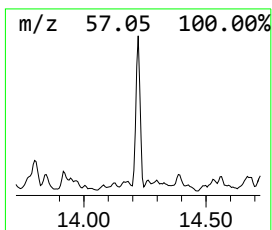
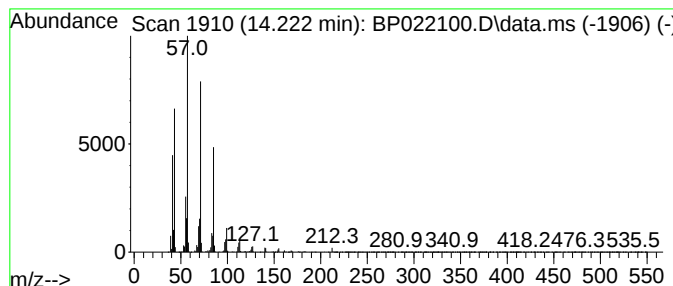
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.222	17.47 ng/ul	1325490	Acenaphthene-d10	14.310	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	90
2	Hexadecane	226	C16H34	000544-76-3	86
3	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	80
4	Decane, 2,3,5,8-tetramethyl-	198	C14H30	192823-15-7	72
5	Nonane, 4-methyl-5-propyl-	184	C13H28	062185-55-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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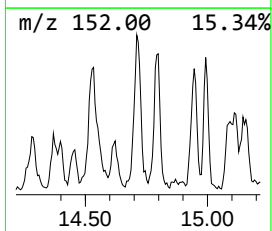
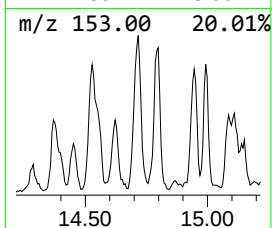
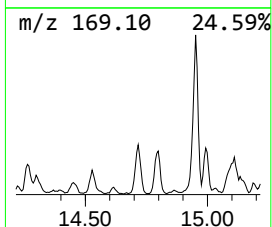
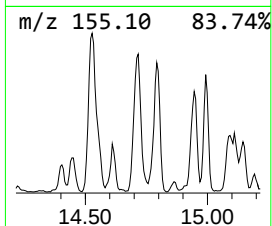
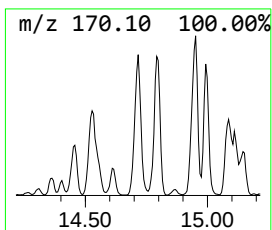
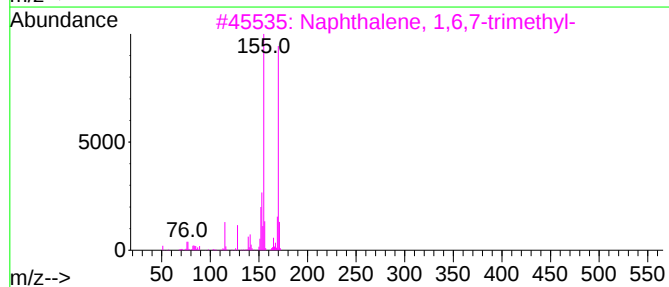
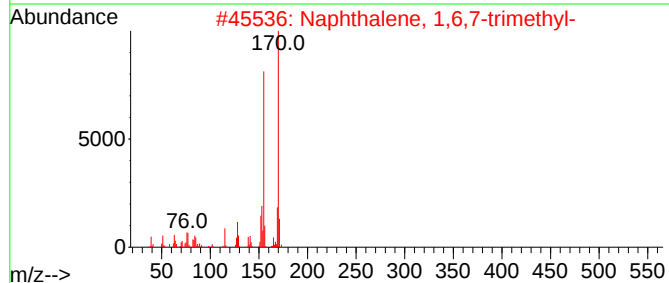
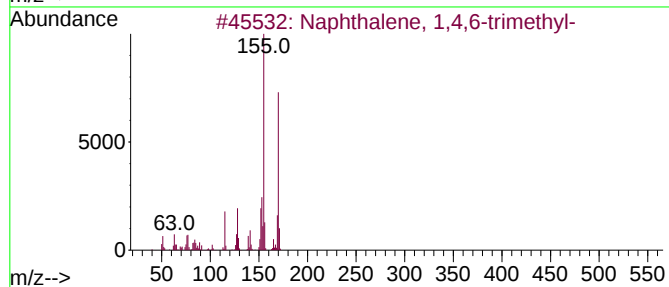
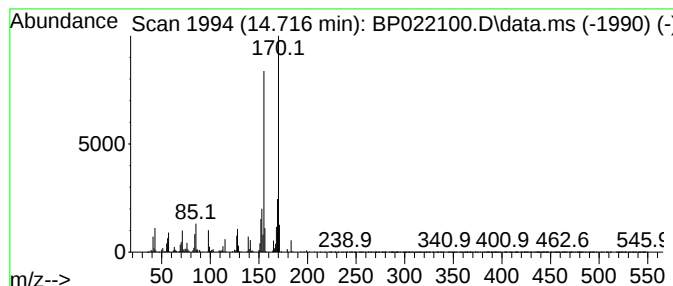
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Quant Title : SVOA CALIBRATION

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

Peak Number 41 Naphthalene, 1,4,6-trimethyl- Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.716	10.65 ng/ul	807844	Acenaphthene-d10	14.310

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	93
2		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
3		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
4		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	93
5		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

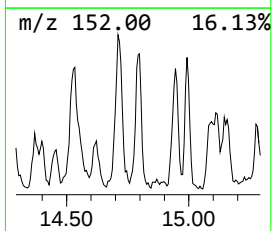
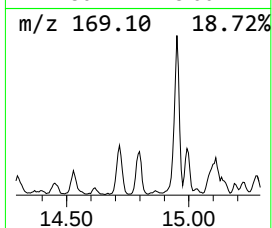
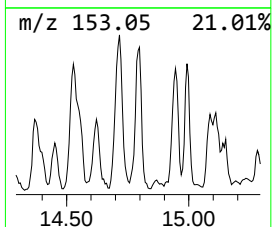
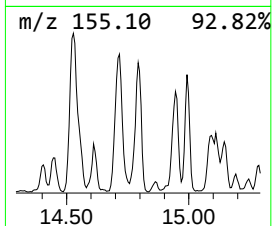
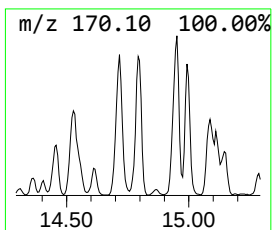
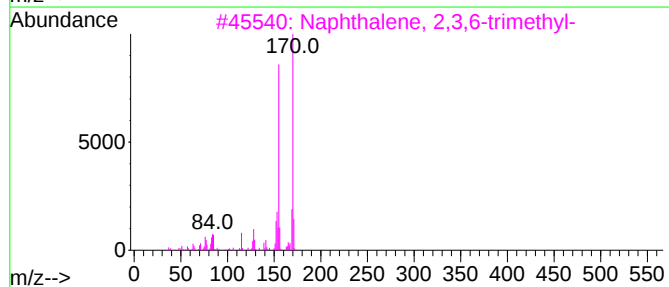
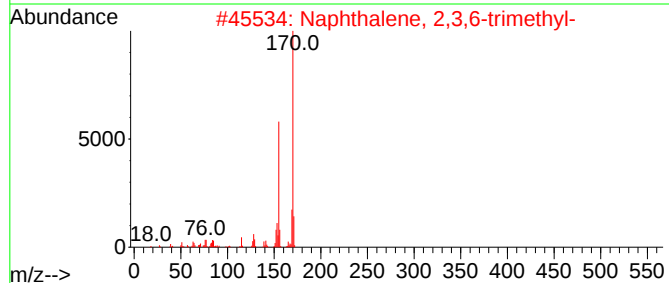
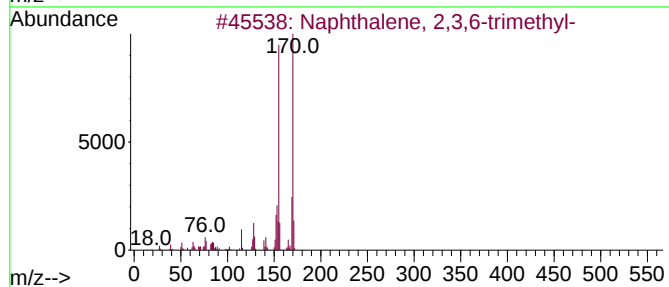
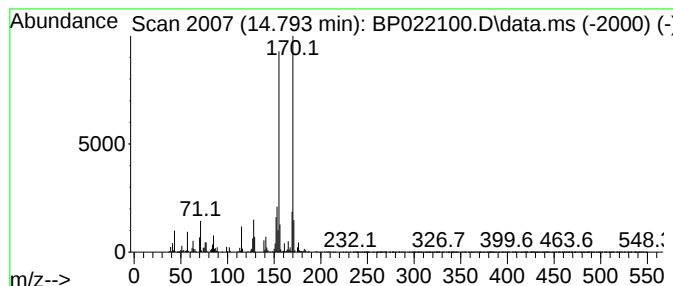
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 42 Naphthalene, 2,3,6-trimethyl- Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.793	8.23 ng/ul	624233	Acenaphthene-d10	14.310	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
2	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
3	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
4	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
5	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC75ME

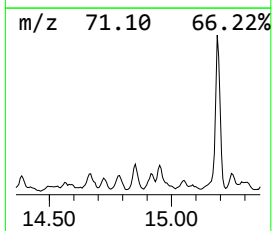
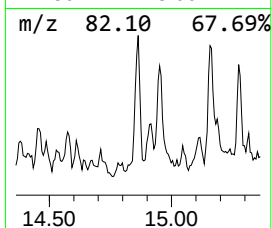
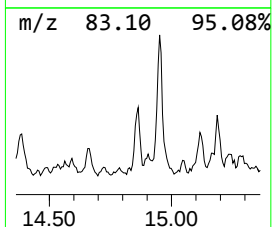
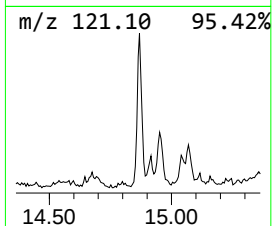
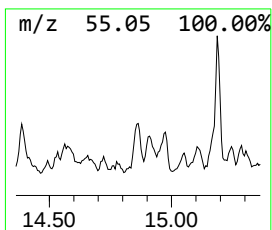
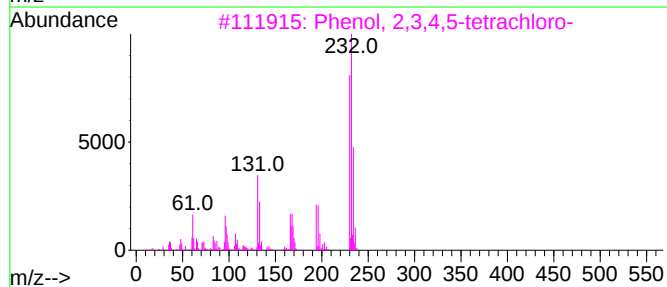
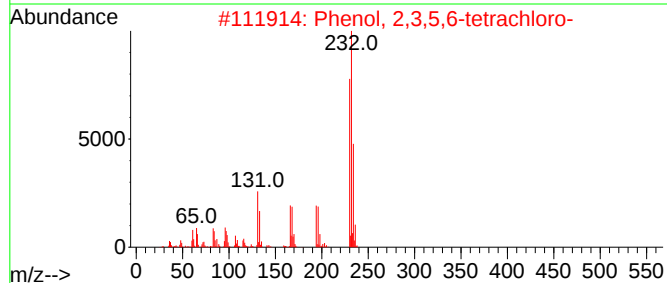
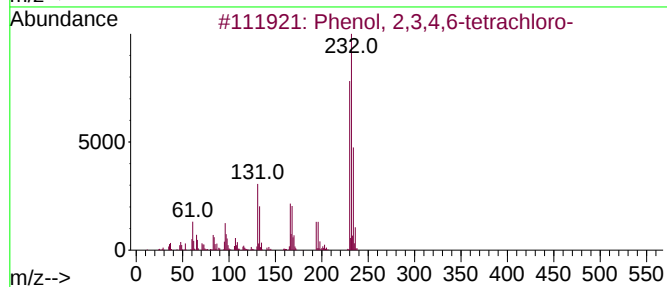
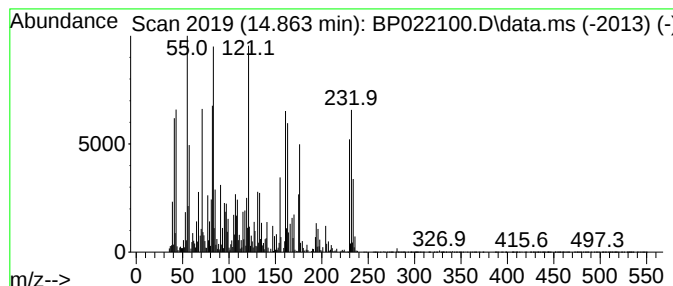
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 Phenol, 2,3,5,6-tetrachloro- Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.863	9.02 ng/ul	684525	Acenaphthene-d10	14.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	92
2			Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	92
3			Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	91
4			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	68
5			Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

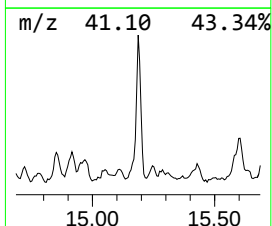
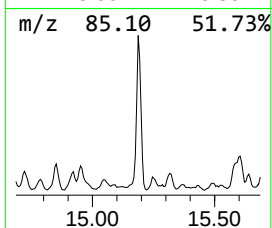
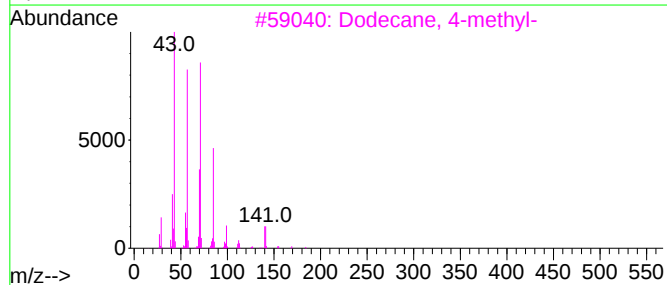
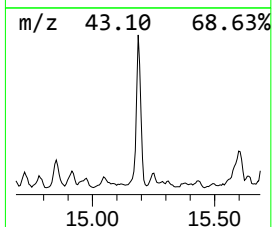
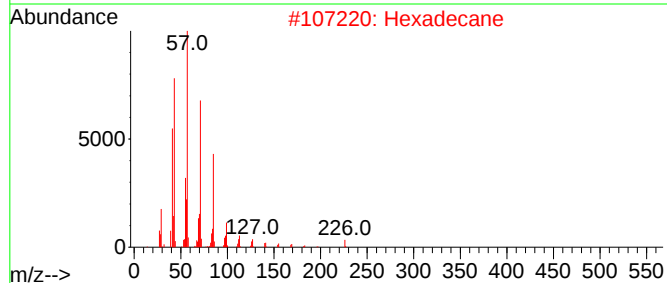
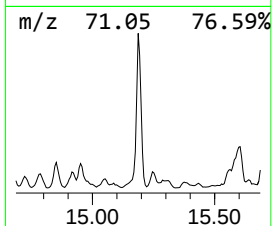
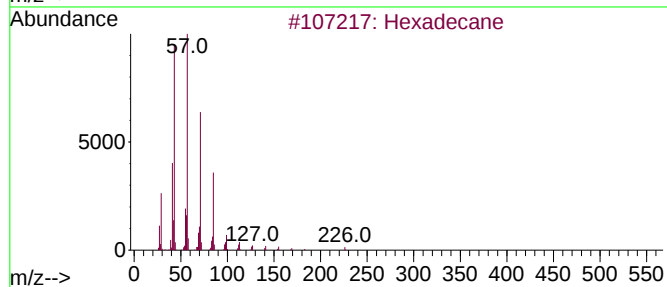
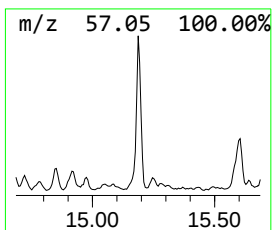
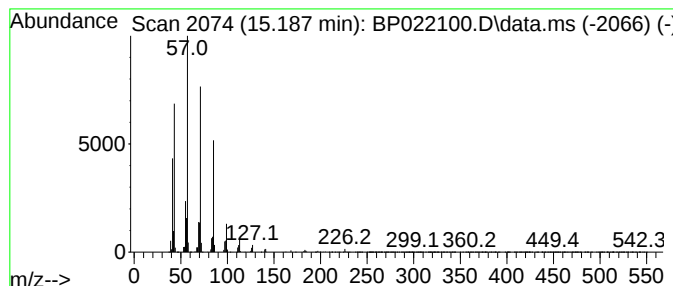
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.187	20.18 ng/ul	1530490	Acenaphthene-d10	14.310	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	94
2	Hexadecane	226	C16H34	000544-76-3	91
3	Dodecane, 4-methyl-	184	C13H28	006117-97-1	90
4	Nonadecane	268	C19H40	000629-92-5	90
5	Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC75ME

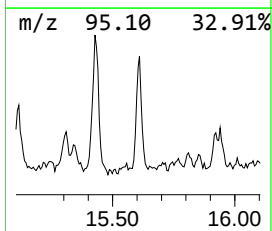
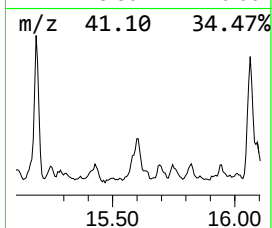
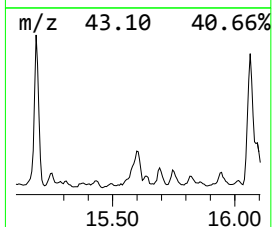
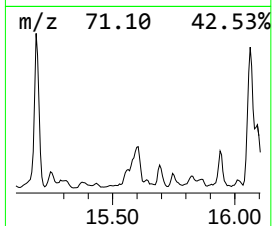
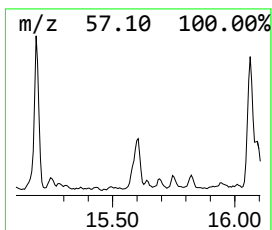
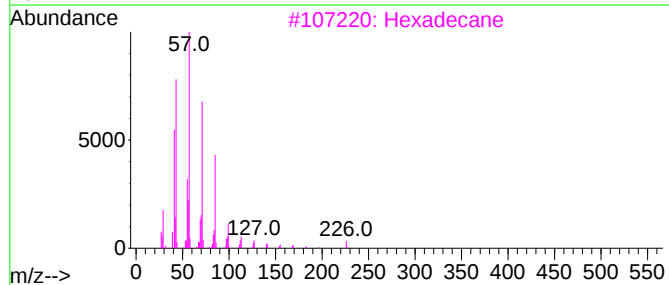
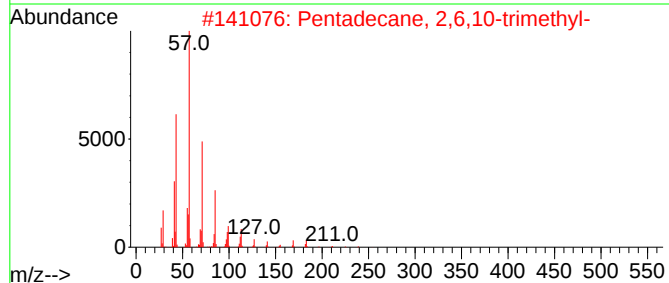
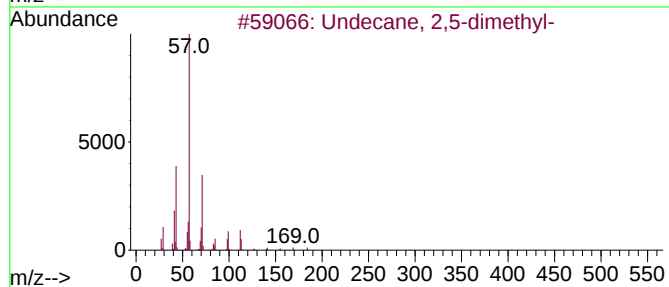
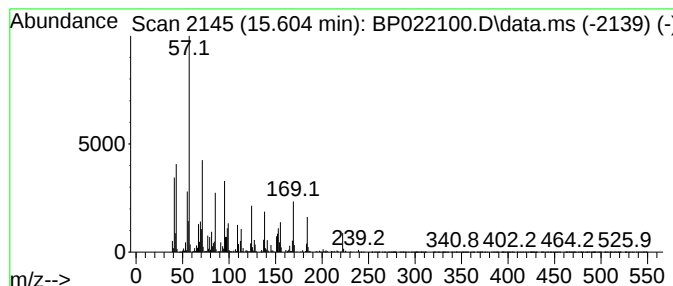
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Quant Title : SVOA CALIBRATION

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

Peak Number 45 (DEL) Alkane: Straight-Chai... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.604	12.83 ng/ul	972861	Acenaphthene-d10	14.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	38
2			Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	38
3			Hexadecane	226	C16H34	000544-76-3	30
4			Hentriacontane	437	C31H64	000630-04-6	30
5			Hexacosane	366	C26H54	000630-01-3	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC75ME

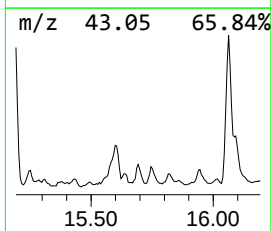
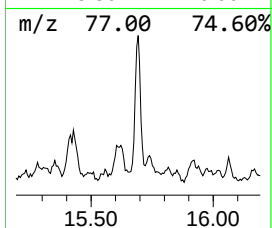
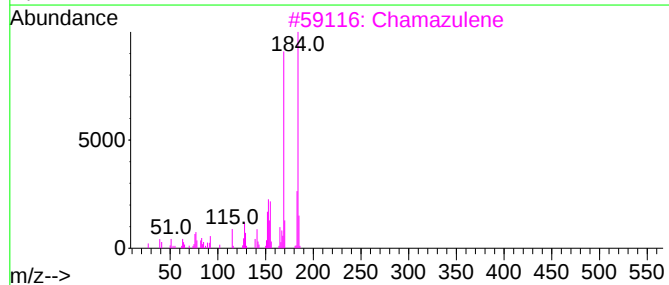
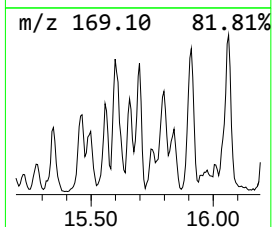
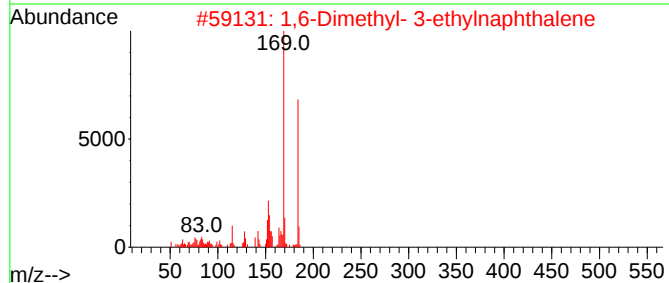
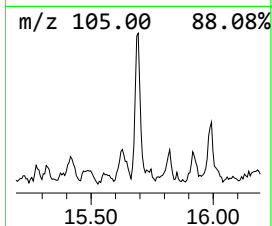
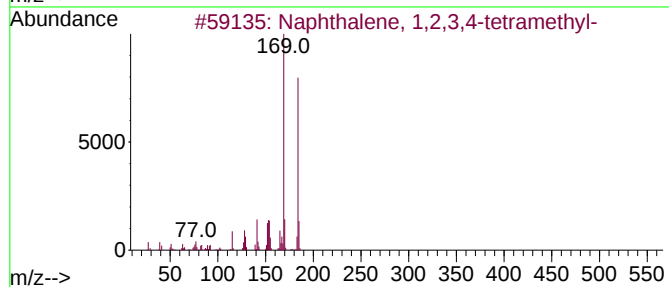
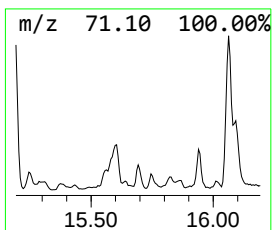
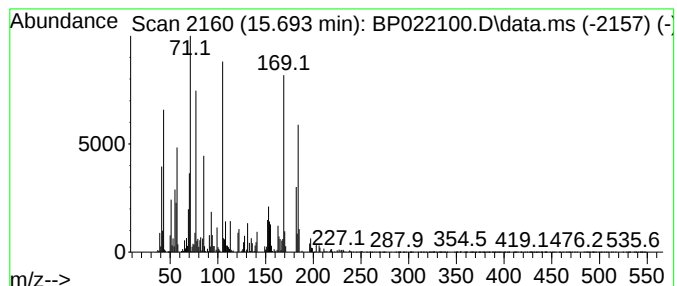
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Quant Title : SVOA CALIBRATION

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

Peak Number 46 Naphthalene, 1,2,3,4-tetram... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.693	7.04 ng/ul	533998	Acenaphthene-d10	14.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetramethyl-	184	C14H16	003031-15-0	70
2			1,6-Dimethyl- 3-ethylnaphthalene	184	C14H16	1000383-71-8	56
3			Chamazulene	184	C14H16	000529-05-5	46
4			3,4,4-Trimethyl-5-methylene-2-et...	184	C9H16N2S	037120-36-8	45
5			1,4,5,8-Tetramethylnaphthalene	184	C14H16	002717-39-7	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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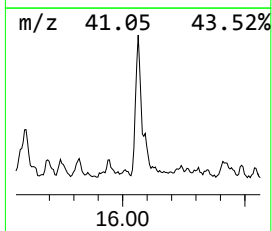
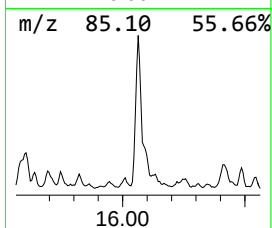
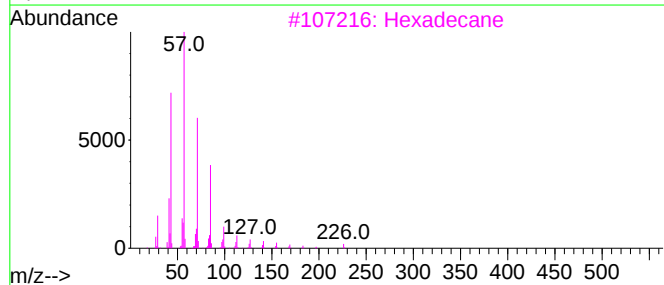
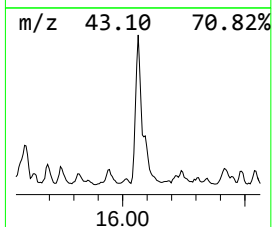
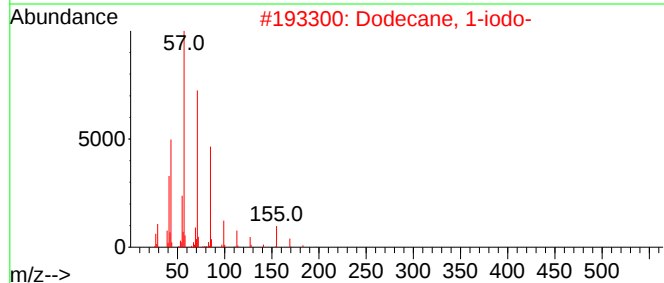
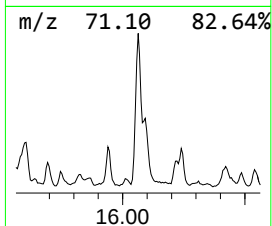
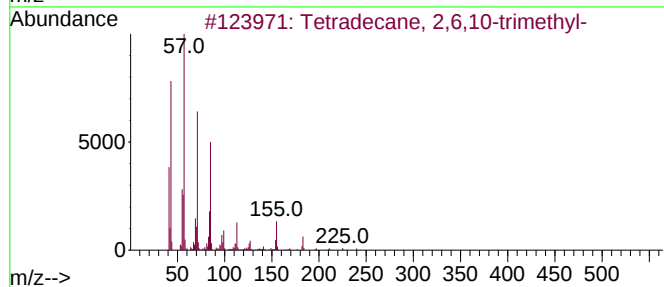
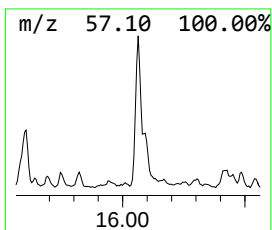
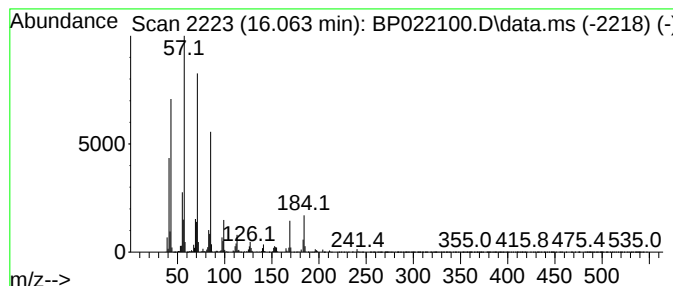
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Quant Title : SVOA CALIBRATION

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

Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.063	21.20 ng/ul	1682060	Phenanthrene-d10	17.110

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	80
2		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80
3		Hexadecane	226	C16H34	000544-76-3	80
4		Dodecane	170	C12H26	000112-40-3	74
5		Tetradecane	198	C14H30	000629-59-4	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

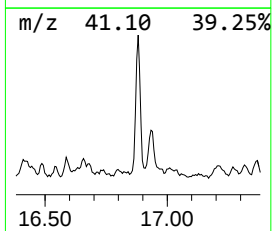
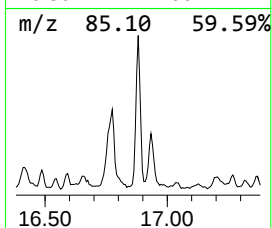
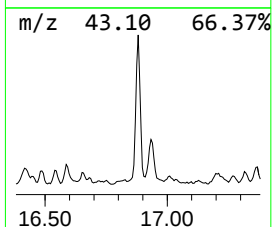
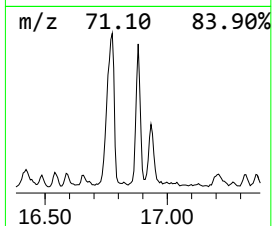
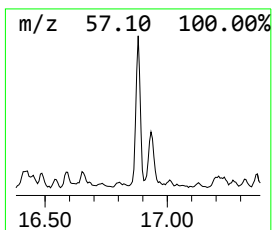
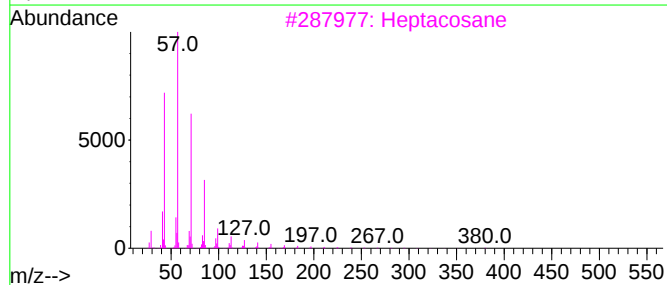
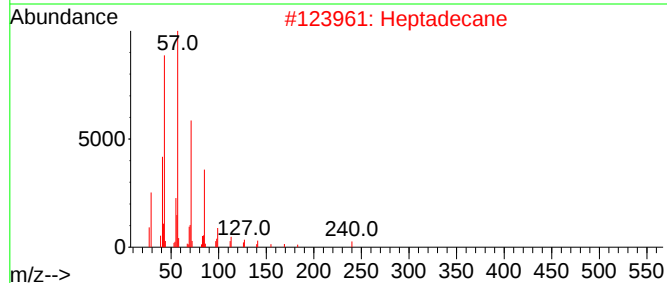
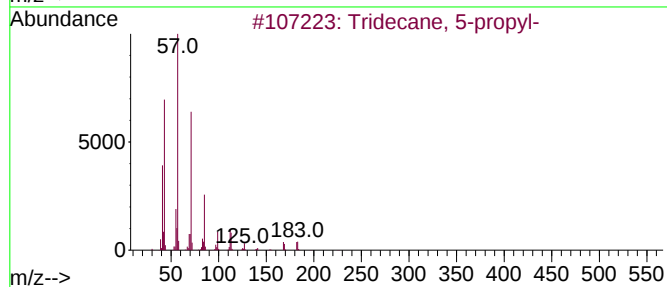
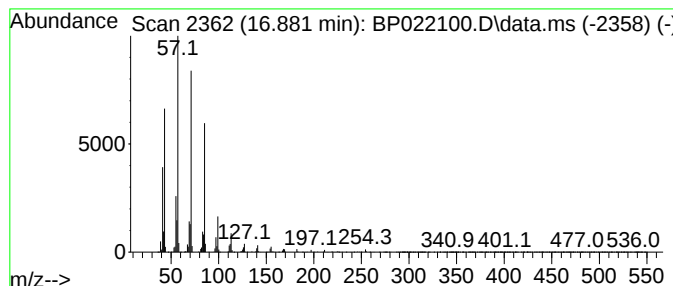
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 48 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.881	15.98 ng/ul	1267830	Phenanthrene-d10		17.110	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tridecane, 5-propyl-		226	C16H34	055045-11-9	86
2	Heptadecane		240	C17H36	000629-78-7	83
3	Heptacosane		380	C27H56	000593-49-7	80
4	Eicosane		282	C20H42	000112-95-8	72
5	Tetradecane		198	C14H30	000629-59-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC75ME

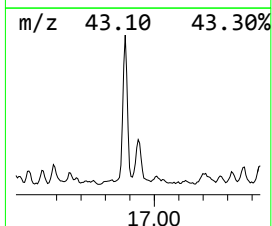
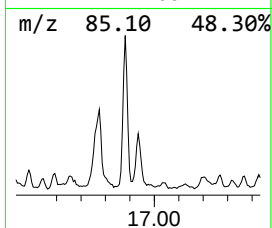
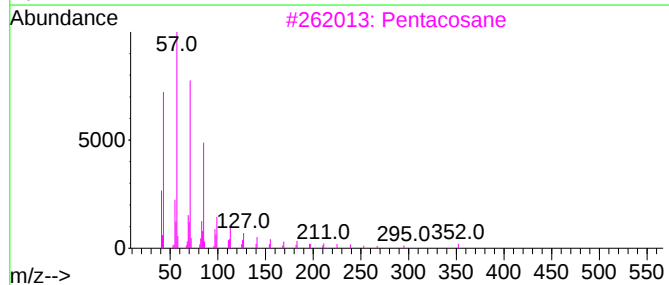
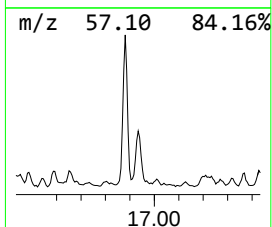
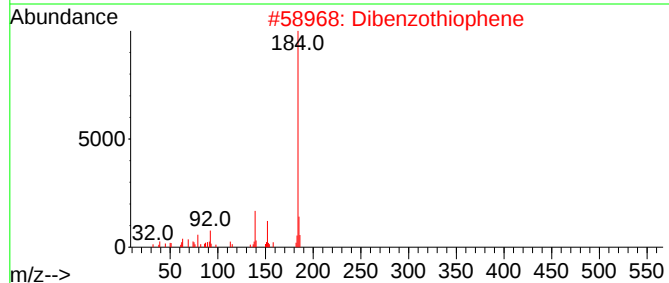
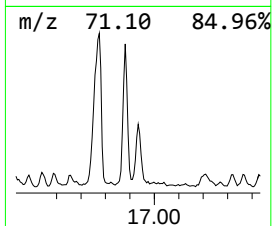
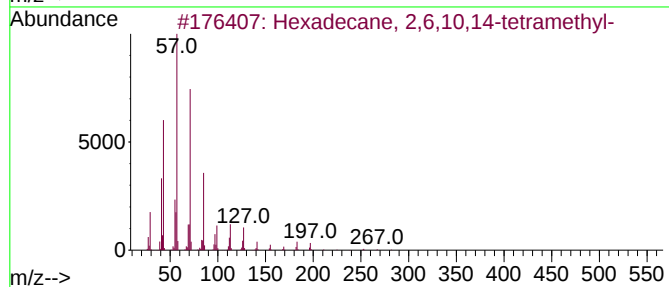
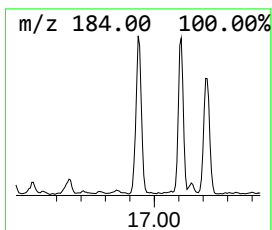
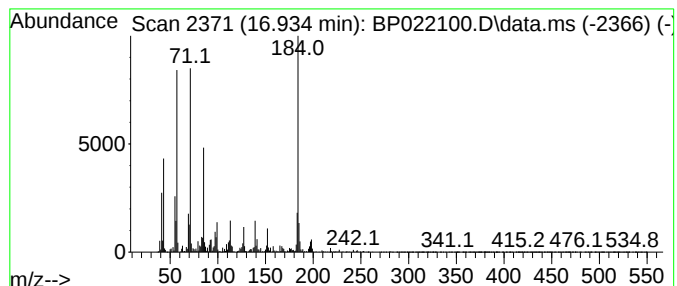
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Quant Title : SVOA CALIBRATION

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

Peak Number 49 (DEL) Alkane: Straight-Chai... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.934	11.89 ng/ul	943159	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	93
2			Dibenzothiophene	184	C12H8S	000132-65-0	62
3			Pentacosane	352	C25H52	000629-99-2	50
4			Dibenzothiophene	184	C12H8S	000132-65-0	46
5			Dibenzothiophene	184	C12H8S	000132-65-0	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

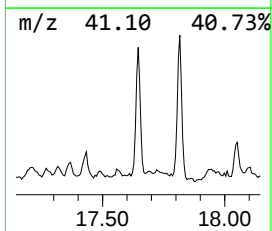
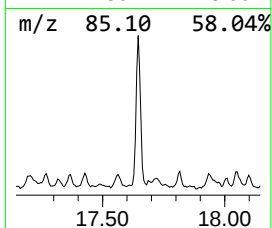
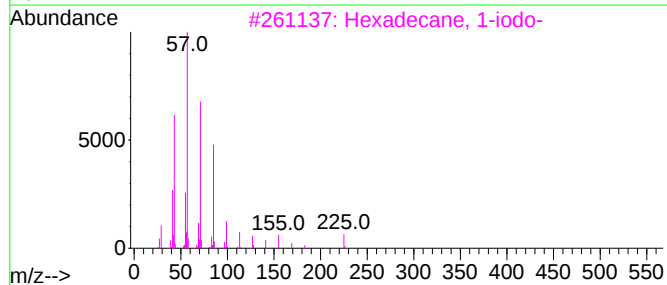
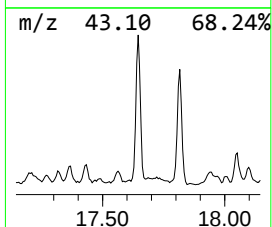
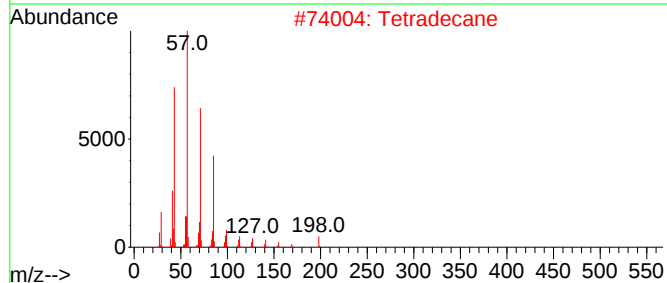
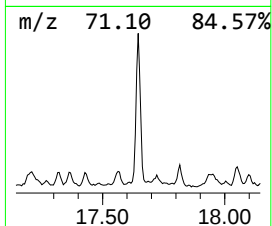
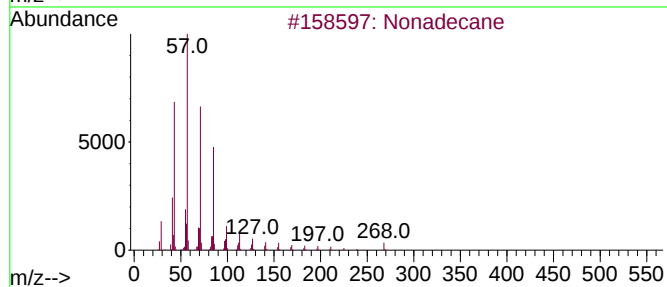
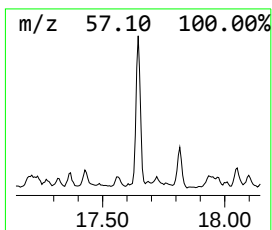
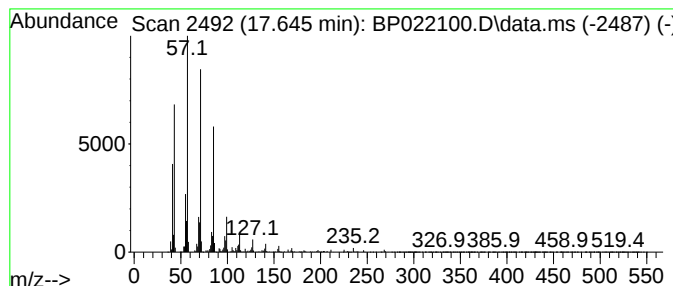
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BNA_P
ClientSampleId :
DDC75ME

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 50 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.645	16.02 ng/ul	1271230	Phenanthrene-d10		17.110	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	91
2	Tetradecane		198	C14H30	000629-59-4	91
3	Hexadecane, 1-iodo-		352	C16H33I	000544-77-4	90
4	Hexadecane		226	C16H34	000544-76-3	90
5	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC75ME

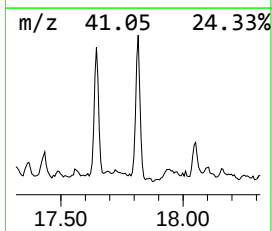
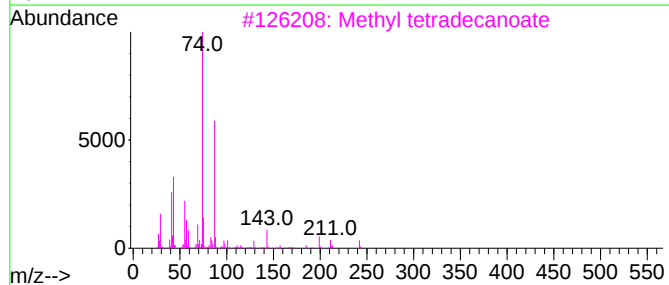
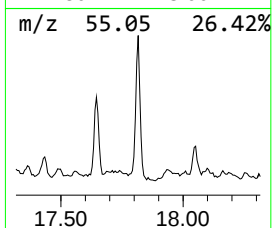
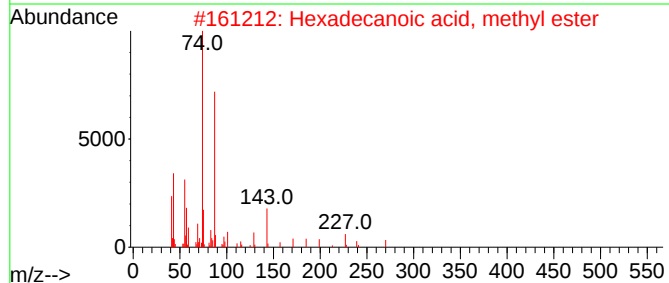
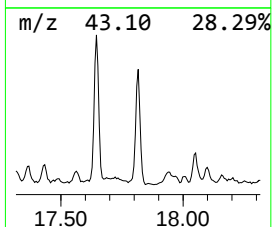
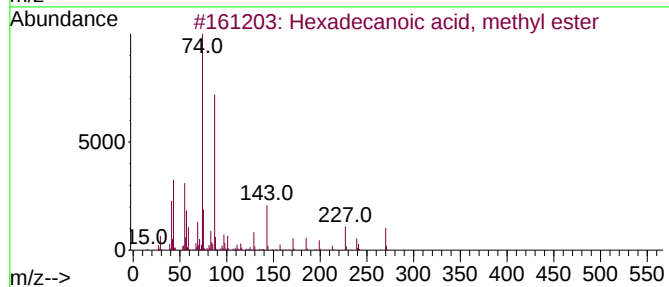
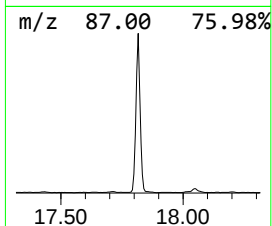
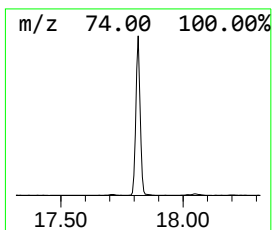
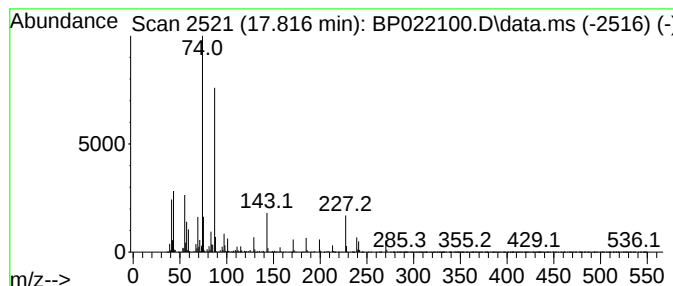
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Quant Title : SVOA CALIBRATION

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

Peak Number 52 Hexadecanoic acid, methyl e... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.816	25.35 ng/ul	2011500	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	97
2			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	96
3			Methyl tetradecanoate	242	C15H30O2	000124-10-7	91
4			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	91
5			Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

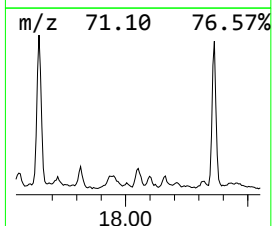
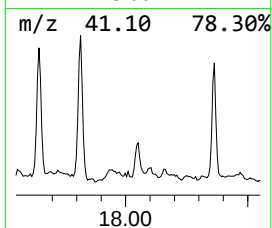
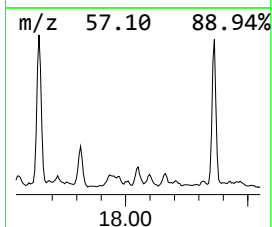
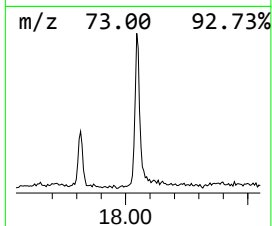
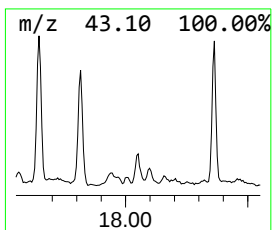
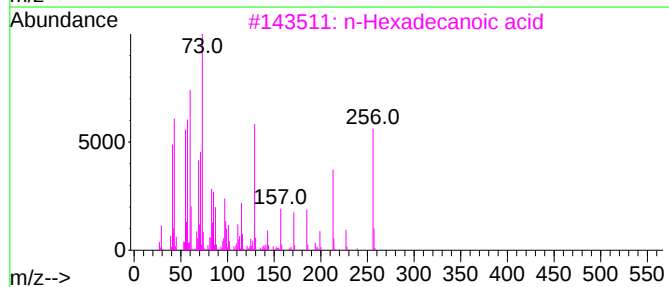
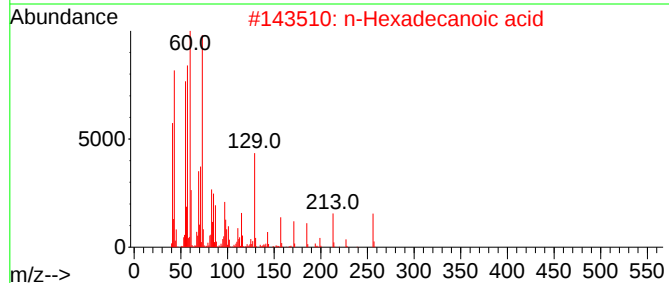
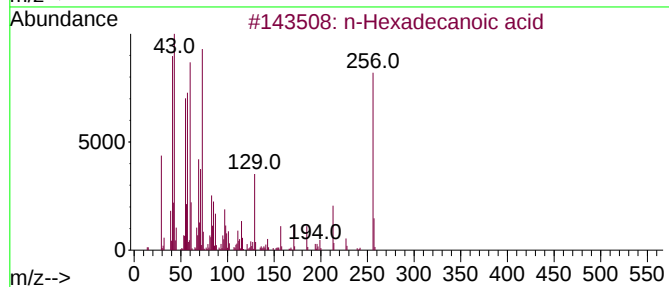
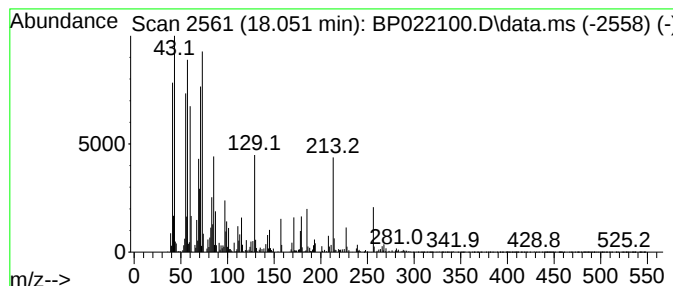
Instrument :
BNA_P
ClientSampleId :
DDC75ME

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 53 n-Hexadecanoic acid Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.			
18.051	8.17 ng/ul	648015	Phenanthrene-d10	17.110			
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	89		
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87		
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	53		
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	46		
5	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	42		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

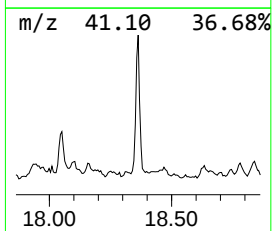
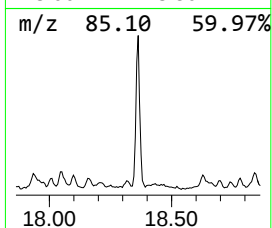
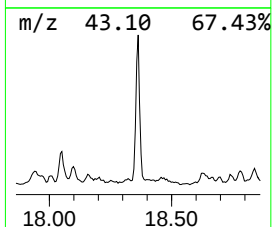
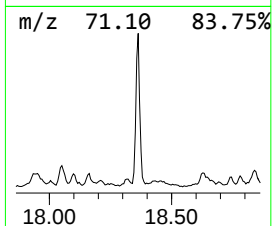
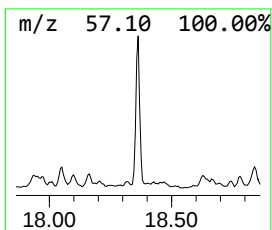
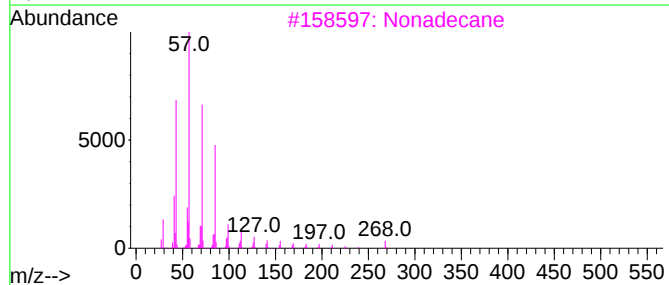
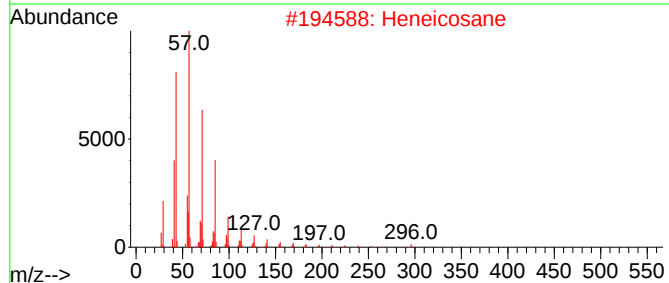
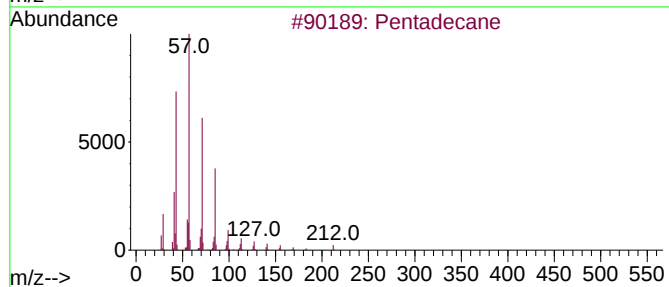
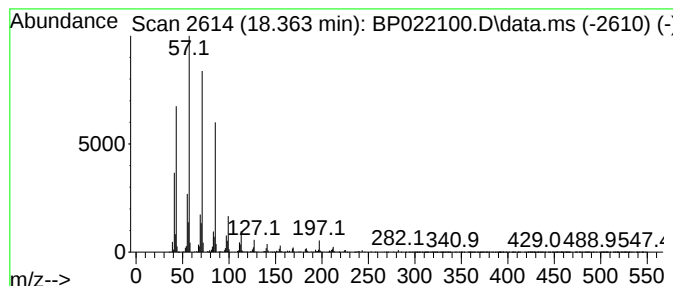
Instrument :
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ClientSampleId :
DDC75ME

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 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.363	14.09 ng/ul	1117610	Phenanthrene-d10		17.110	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecane		212	C15H32	000629-62-9	97
2	Heneicosane		296	C21H44	000629-94-7	91
3	Nonadecane		268	C19H40	000629-92-5	90
4	Hexadecane, 1-iodo-		352	C16H33I	000544-77-4	90
5	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

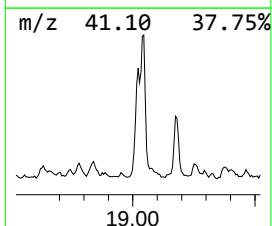
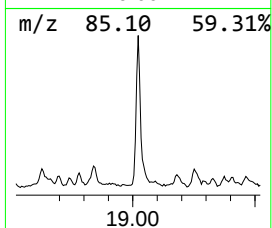
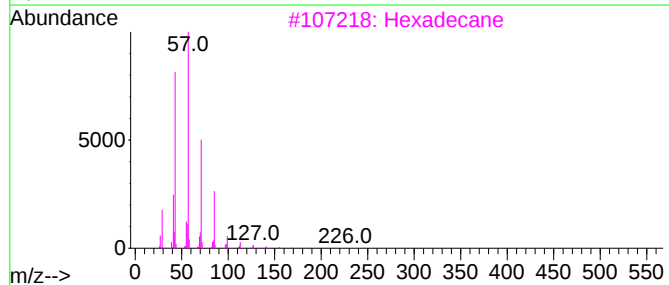
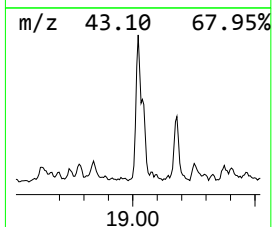
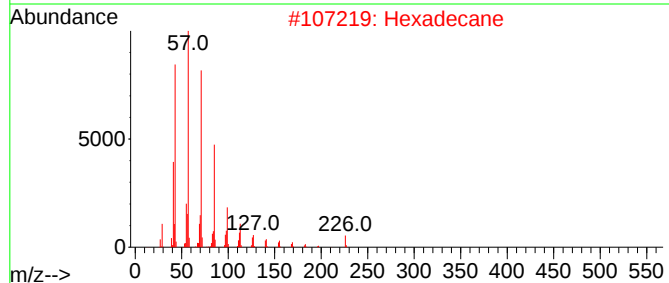
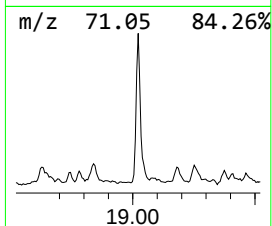
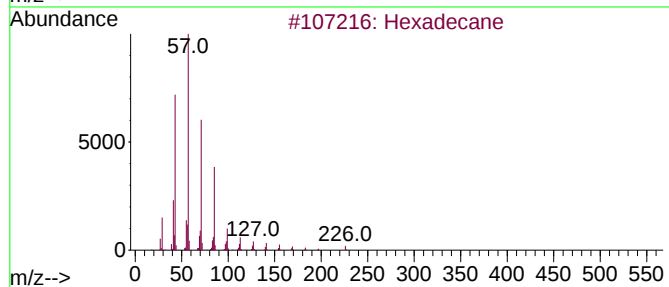
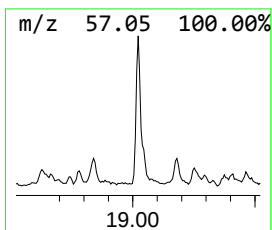
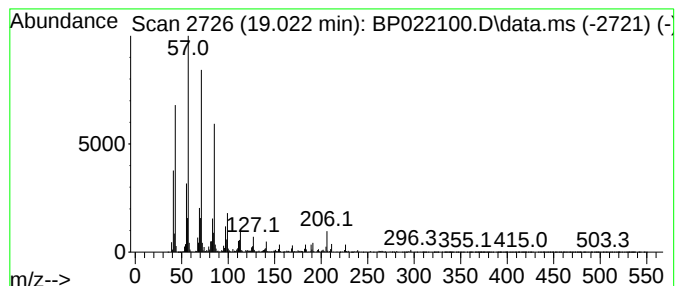
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 55 (DEL) Alkane: Straight-Chai... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.022	14.27 ng/ul	1132020	Phenanthrene-d10		17.110	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	93
2	Hexadecane		226	C16H34	000544-76-3	93
3	Hexadecane		226	C16H34	000544-76-3	92
4	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	91
5	Hexacosane		366	C26H54	000630-01-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC75ME

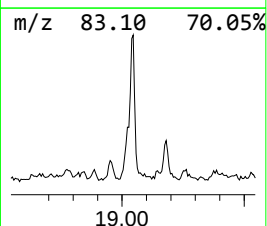
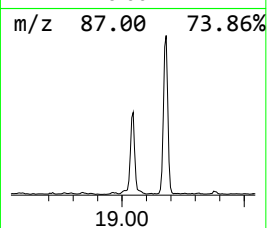
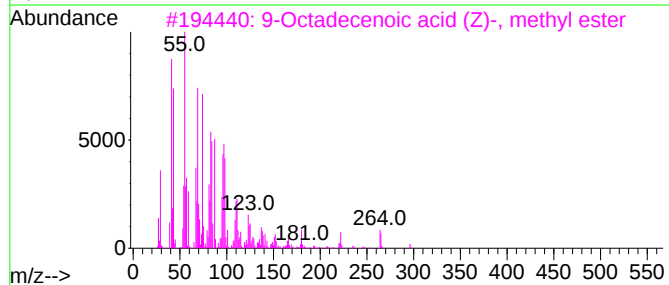
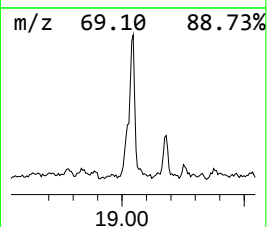
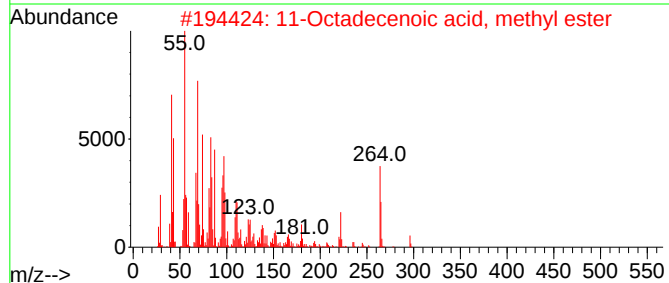
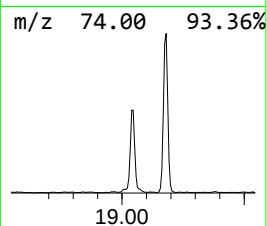
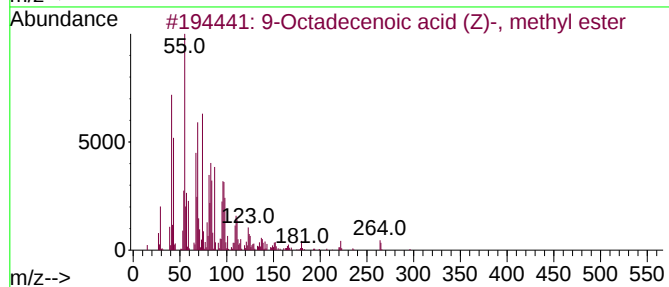
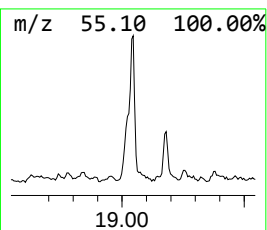
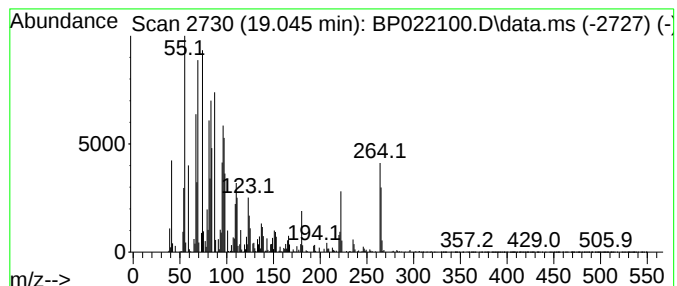
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Quant Title : SVOA CALIBRATION

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

Peak Number 56 9-Octadecenoic acid (Z)-, m... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.045	19.49 ng/ul	1546130	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	93
2			11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	87
3			9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	83
4			8-Octadecenoic acid, methyl ester	296	C19H36O2	002345-29-1	76
5			10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

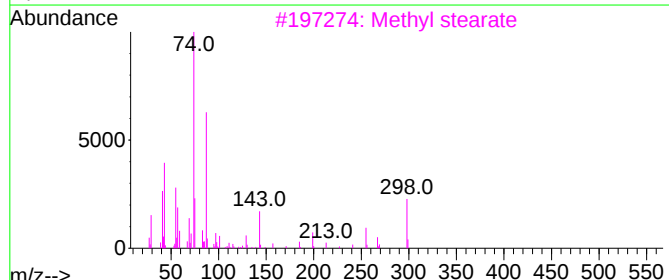
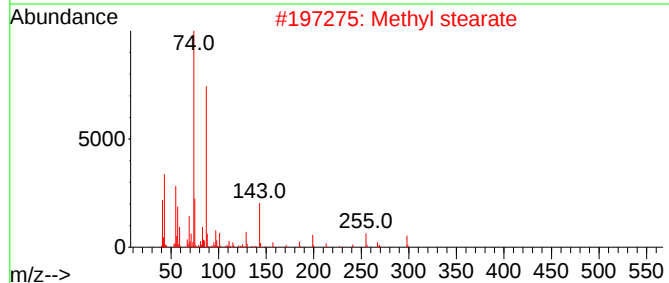
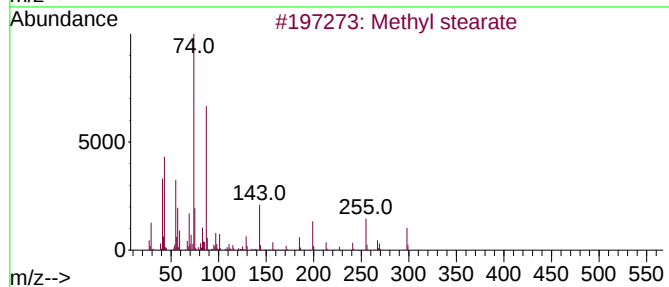
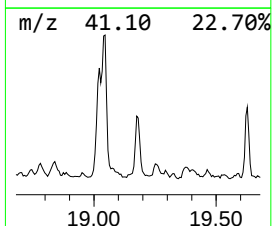
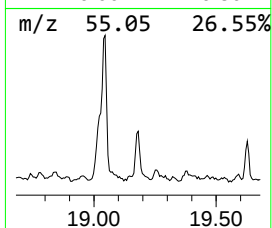
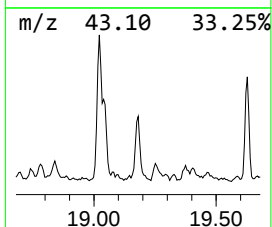
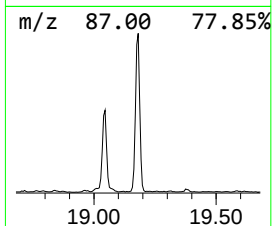
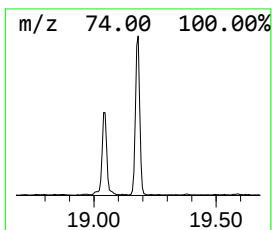
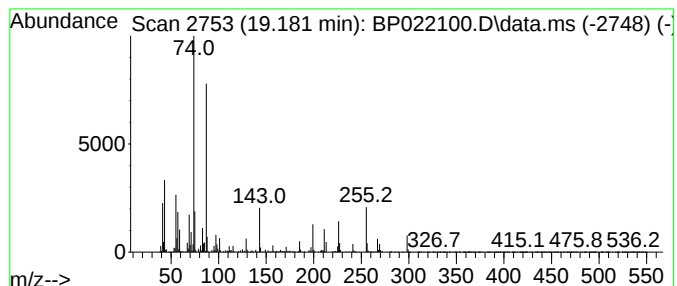
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ClientSampleId :
DDC75ME

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 Methyl stearate Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.181	12.92 ng/ul	1025330	Phenanthrene-d10		17.110	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Methyl stearate		298	C19H38O2	000112-61-8	99
2	Methyl stearate		298	C19H38O2	000112-61-8	97
3	Methyl stearate		298	C19H38O2	000112-61-8	94
4	Heptadecanoic acid, 16-methyl-, ...		298	C19H38O2	005129-61-3	91
5	Methyl stearate		298	C19H38O2	000112-61-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC75ME

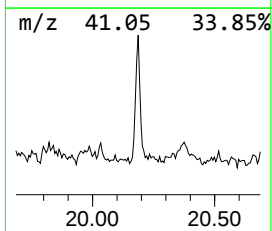
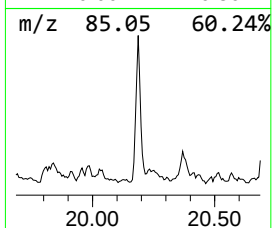
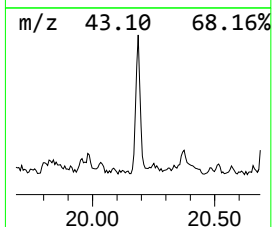
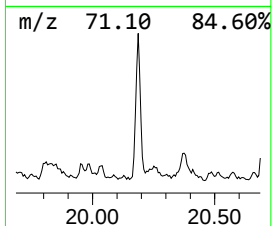
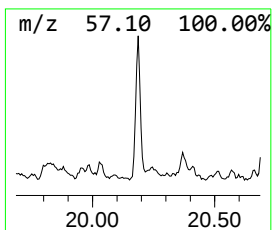
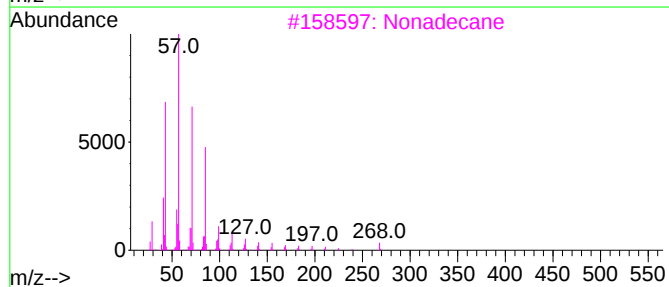
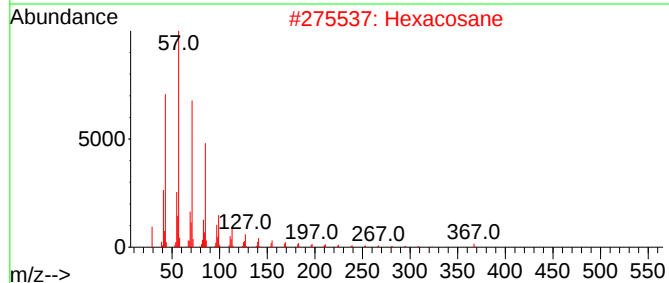
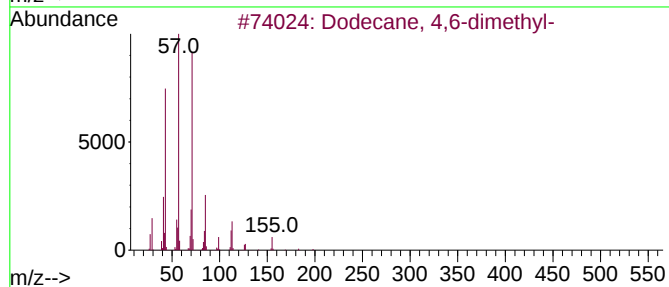
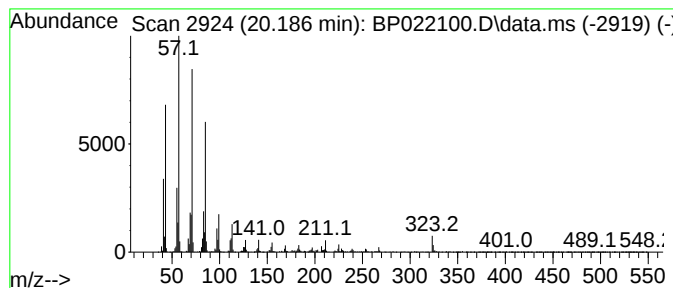
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Quant Title : SVOA CALIBRATION

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

Peak Number 58 (DEL) Alkane: Straight-Chai... Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.186	6.82 ng/ul	693249	Chrysene-d12	21.539

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	93
2		Hexacosane	366	C26H54	000630-01-3	91
3		Nonadecane	268	C19H40	000629-92-5	91
4		Triacontane	422	C30H62	000638-68-6	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC75ME

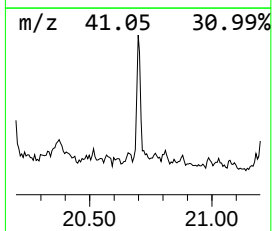
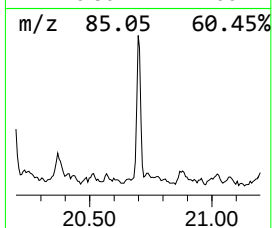
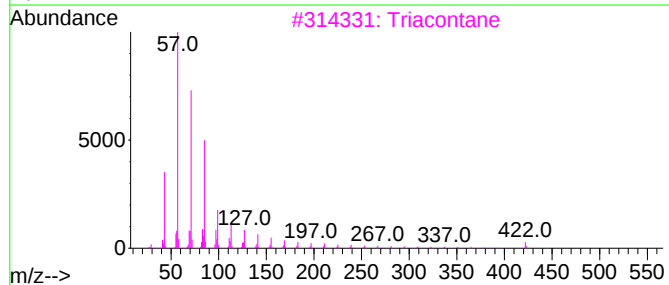
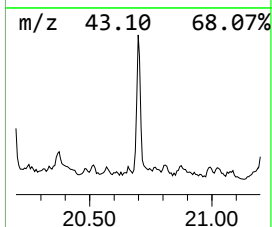
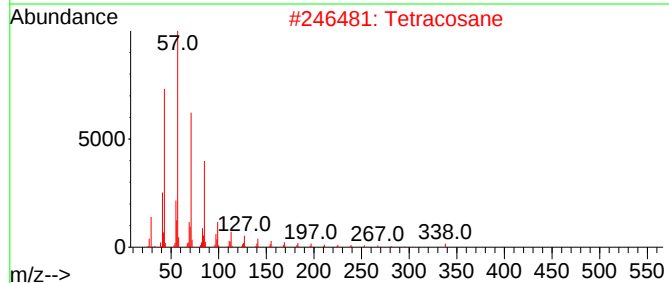
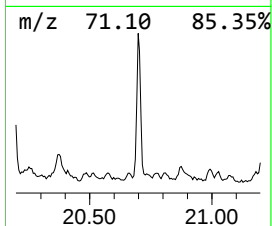
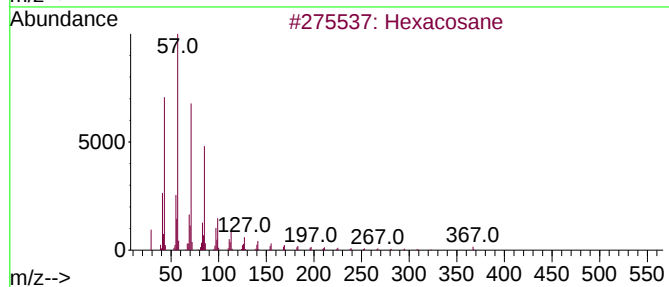
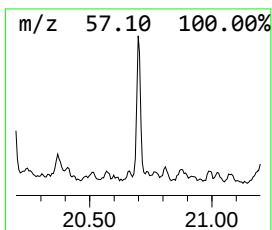
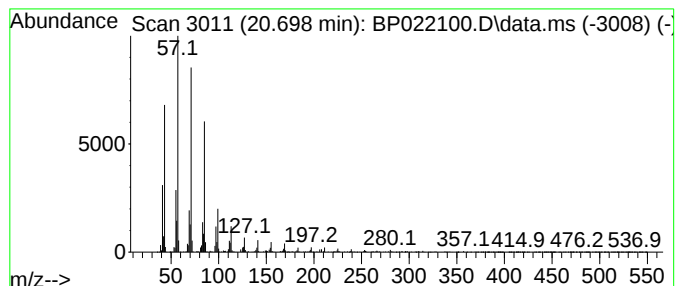
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Quant Title : SVOA CALIBRATION

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

Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.698	5.44 ng/ul	553054	Chrysene-d12	21.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	94
2			Tetracosane	338	C24H50	000646-31-1	91
3			Triacontane	422	C30H62	000638-68-6	91
4			Heneicosane	296	C21H44	000629-94-7	91
5			Docosane, 11-butyl-	366	C26H54	013475-76-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC75ME

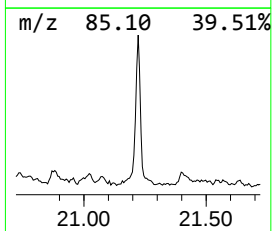
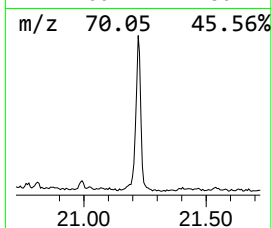
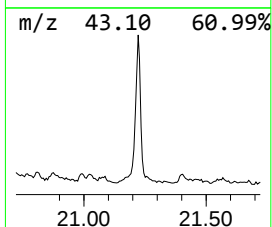
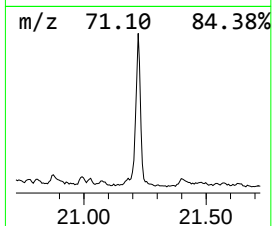
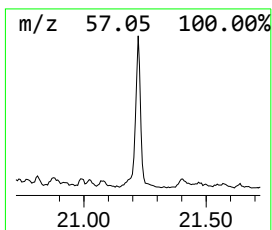
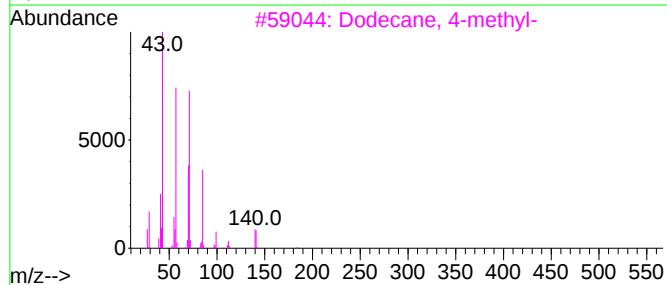
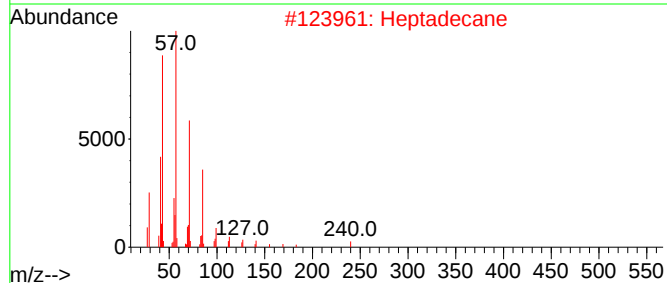
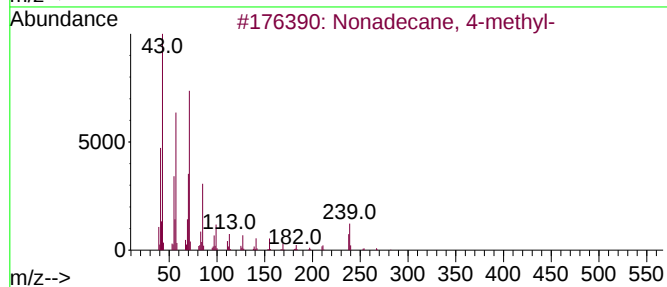
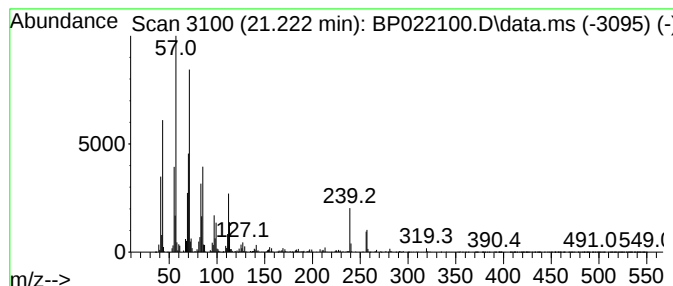
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Quant Title : SVOA CALIBRATION

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

Peak Number 60 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.222	13.87 ng/ul	1410160	Chrysene-d12	21.539

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane, 4-methyl-	282	C20H42	025117-27-5	90
2		Heptadecane	240	C17H36	000629-78-7	60
3		Dodecane, 4-methyl-	184	C13H28	006117-97-1	58
4		Heptadecane, 7-methyl-	254	C18H38	020959-33-5	58
5		Carbonic acid, dodecyl 2-ethylhe...	342	C21H42O3	1000383-13-9	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC75ME

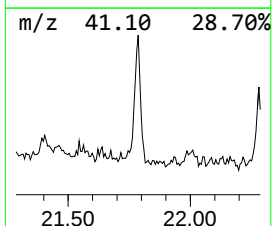
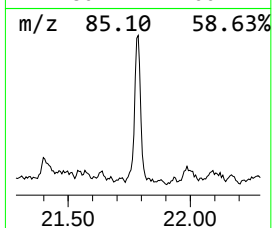
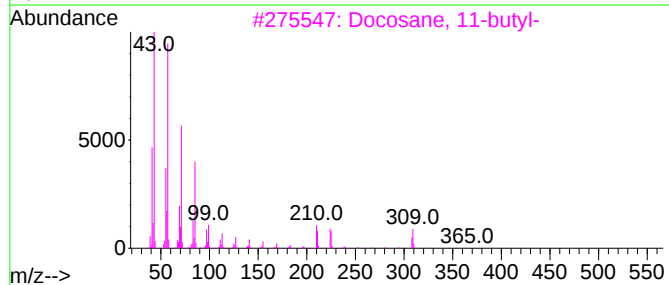
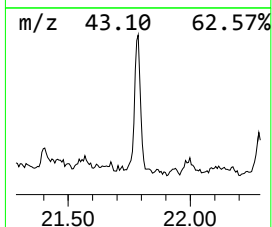
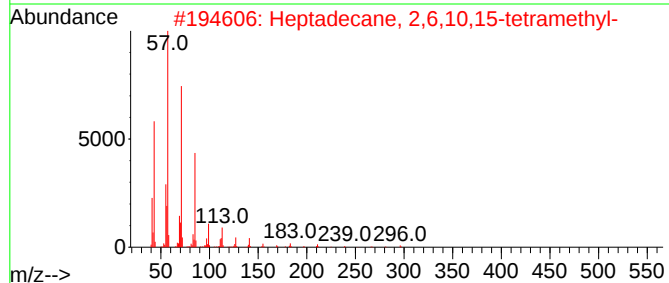
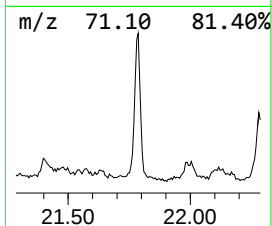
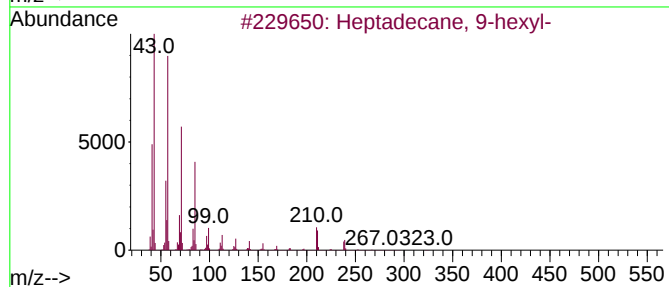
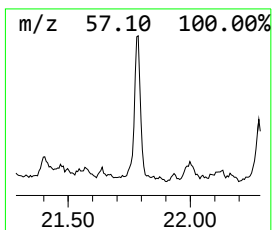
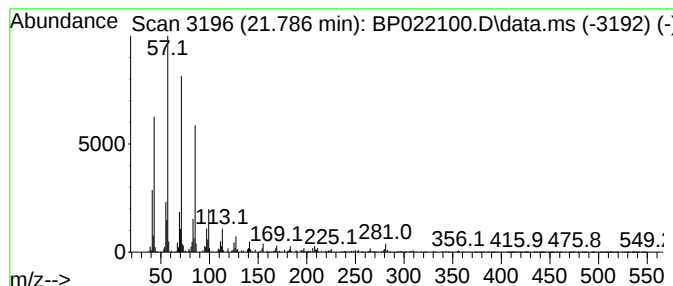
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Quant Title : SVOA CALIBRATION

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

Peak Number 61 (DEL) Alkane: Straight-Chai... Concentration Rank 56

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.786	5.12 ng/ul	520725	Chrysene-d12	21.539

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	91
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	91
4		Hexacosane	366	C26H54	000630-01-3	90
5		Octadecane, 2-methyl-	268	C19H40	001560-88-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022100.D
Acq On : 28 Sep 2024 11:42
Operator : RC/JU
Sample : P3910-02ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

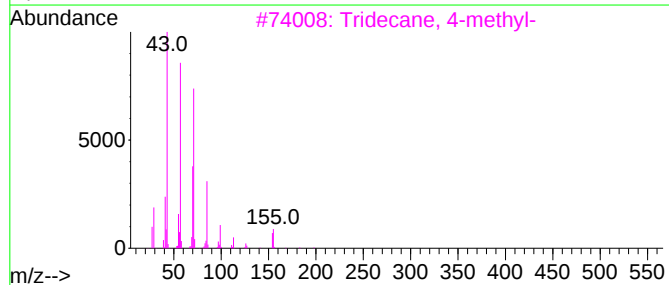
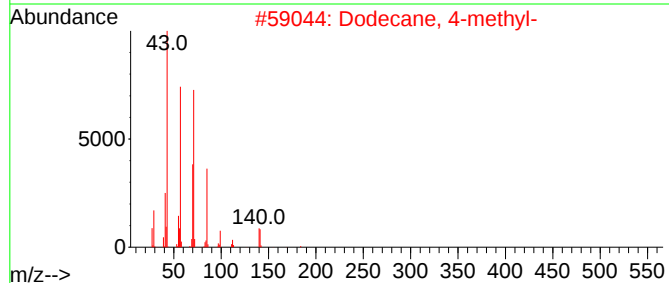
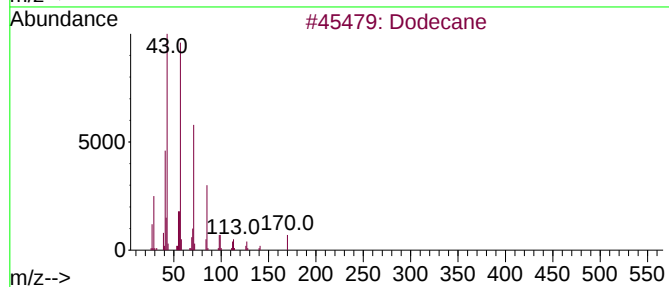
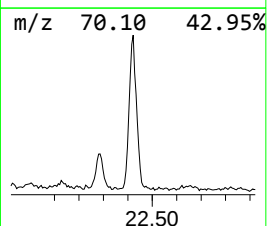
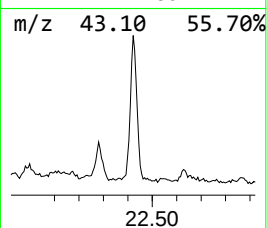
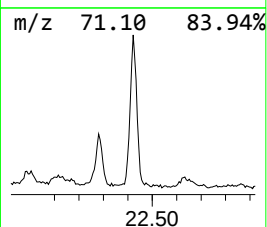
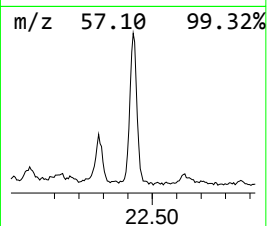
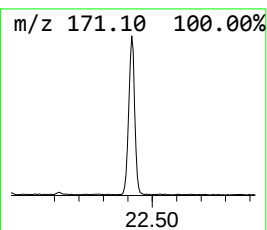
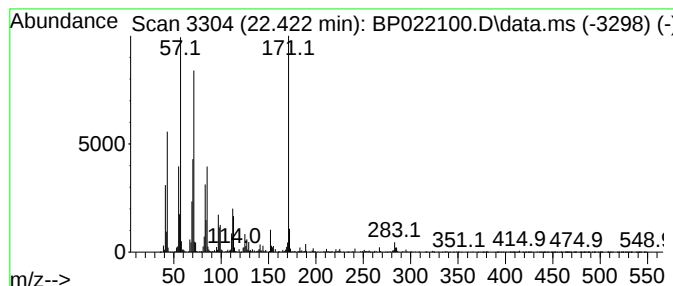
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 62 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.422	10.74 ng/ul	1092440	Chrysene-d12	21.539	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane	170	C12H26	000112-40-3	50
2	Dodecane, 4-methyl-	184	C13H28	006117-97-1	49
3	Tridecane, 4-methyl-	198	C14H30	026730-12-1	46
4	Dodecane	170	C12H26	000112-40-3	45
5	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
 Data File : BP022100.D
 Acq On : 28 Sep 2024 11:42
 Operator : RC/JU
 Sample : P3910-02ME
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DDC75ME

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...	5.758	17.0	ng/ul	1088830	1	7.699	1280030	20.0
(DEL) Alkane: C...	6.369	8.5	ng/ul	542179	1	7.699	1280030	20.0
Benzene, propyl-	6.711	19.3	ng/ul	1234260	1	7.699	1280030	20.0
1-Iodo-2-methyl...	6.769	12.3	ng/ul	787313	1	7.699	1280030	20.0
Benzene, 1-ethy...	6.811	22.9	ng/ul	1464570	1	7.699	1280030	20.0
Benzene, 1,2,3-...	6.940	11.2	ng/ul	713502	1	7.699	1280030	20.0
Benzene, 1-ethy...	7.105	13.7	ng/ul	875462	1	7.699	1280030	20.0
2-Heptanone, 4,...	7.140	17.1	ng/ul	1095930	1	7.699	1280030	20.0
(DEL) Alkane: S...	7.340	91.0	ng/ul	5825620	1	7.699	1280030	20.0
(DEL) Alkane: S...	7.628	8.8	ng/ul	561012	1	7.699	1280030	20.0
2-Ethylhexyl hy...	7.769	12.0	ng/ul	767510	1	7.699	1280030	20.0
Benzene, 1,2,4-...	7.816	13.1	ng/ul	838166	1	7.699	1280030	20.0
(DEL) Alkane: C...	7.993	8.3	ng/ul	532355	1	7.699	1280030	20.0
Cyclohexanone, ...	8.128	33.3	ng/ul	2129730	1	7.699	1280030	20.0
unknown-01	8.175	7.3	ng/ul	466535	1	7.699	1280030	20.0
Benzene, 1-meth...	8.234	16.1	ng/ul	1030400	1	7.699	1280030	20.0
Benzene, 1,4-di...	8.334	19.0	ng/ul	1218220	1	7.699	1280030	20.0
(DEL) Alkane: S...	8.458	14.3	ng/ul	915367	1	7.699	1280030	20.0
Benzene, 1-meth...	8.493	11.9	ng/ul	761872	1	7.699	1280030	20.0
Benzene, 4-ethy...	8.640	9.7	ng/ul	620905	1	7.699	1280030	20.0
o-Cymene	8.687	15.5	ng/ul	994802	1	7.699	1280030	20.0
(DEL) Alkane: S...	8.916	63.4	ng/ul	4059680	1	7.699	1280030	20.0
4-Methylphenyl ...	9.034	14.7	ng/ul	939770	1	7.699	1280030	20.0
(DEL) Alkane: C...	11.369	2.1	ng/ul	795126	2	10.458	7431060	20.0
(DEL) Alkane: S...	11.722	2.3	ng/ul	851214	2	10.458	7431060	20.0
unknown-02	12.557	7.4	ng/ul	561272	3	14.310	1517050	20.0
unknown-03	12.881	6.6	ng/ul	498566	3	14.310	1517050	20.0
(DEL) Alkane: S...	13.134	21.4	ng/ul	1621930	3	14.310	1517050	20.0
unknown-04	13.357	15.5	ng/ul	1173810	3	14.310	1517050	20.0
Naphthalene, 2,...	13.487	21.7	ng/ul	1647430	3	14.310	1517050	20.0
Naphthalene, 1,...	13.628	16.4	ng/ul	1244530	3	14.310	1517050	20.0
(DEL) Alkane: S...	13.799	8.4	ng/ul	638868	3	14.310	1517050	20.0
Naphthalene, 1,...	13.863	9.2	ng/ul	694931	3	14.310	1517050	20.0
(DEL) Alkane: S...	14.222	17.5	ng/ul	1325490	3	14.310	1517050	20.0
Naphthalene, 1,...	14.716	10.7	ng/ul	807844	3	14.310	1517050	20.0
Naphthalene, 2,...	14.793	8.2	ng/ul	624233	3	14.310	1517050	20.0
Phenol, 2,3,5,6...	14.863	9.0	ng/ul	684525	3	14.310	1517050	20.0
(DEL) Alkane: S...	15.187	20.2	ng/ul	1530490	3	14.310	1517050	20.0
(DEL) Alkane: S...	15.604	12.8	ng/ul	972861	3	14.310	1517050	20.0
Naphthalene, 1,...	15.693	7.0	ng/ul	533998	3	14.310	1517050	20.0
(DEL) Alkane: S...	16.063	21.2	ng/ul	1682060	4	17.110	1586790	20.0
(DEL) Alkane: S...	16.881	16.0	ng/ul	1267830	4	17.110	1586790	20.0
(DEL) Alkane: S...	16.934	11.9	ng/ul	943159	4	17.110	1586790	20.0
(DEL) Alkane: S...	17.645	16.0	ng/ul	1271230	4	17.110	1586790	20.0
Hexadecanoic ac...	17.816	25.4	ng/ul	2011500	4	17.110	1586790	20.0
n-Hexadecanoic ...	18.051	8.2	ng/ul	648015	4	17.110	1586790	20.0
(DEL) Alkane: S...	18.363	14.1	ng/ul	1117610	4	17.110	1586790	20.0
(DEL) Alkane: S...	19.022	14.3	ng/ul	1132020	4	17.110	1586790	20.0
9-Octadecenoic ...	19.045	19.5	ng/ul	1546130	4	17.110	1586790	20.0
Methyl stearate	19.181	12.9	ng/ul	1025330	4	17.110	1586790	20.0
(DEL) Alkane: S...	20.186	6.8	ng/ul	693249	5	21.539	2033590	20.0
(DEL) Alkane: S...	20.698	5.4	ng/ul	553054	5	21.539	2033590	20.0

(DEL) Alkane: S...	21.222	13.9	ng/ul	1410160	5	21.539	2033590	20.0
(DEL) Alkane: S...	21.786	5.1	ng/ul	520725	5	21.539	2033590	20.0
(DEL) Alkane: S...	22.422	10.7	ng/ul	1092440	5	21.539	2033590	20.0

Instrument :
BNA_P
ClientSampleId :
DDC75ME