

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
 Data File : BP022092.D
 Acq On : 28 Sep 2024 05:10
 Operator : RC/JU
 Sample : P3910-17 10X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DDC63

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: BP022092.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.870	313	320	329	rVB	89538	147328	0.75%	0.224%
2	5.758	464	471	478	rVB2	370262	558495	2.86%	0.849%
3	6.058	517	522	527	rVV	253953	394480	2.02%	0.600%
4	6.217	544	549	556	rBV4	67126	136498	0.70%	0.208%
5	6.287	556	561	566	rVV	125314	178818	0.92%	0.272%
6	6.369	570	575	578	rVV2	95433	197928	1.01%	0.301%
7	6.711	629	633	638	rVV2	177251	286505	1.47%	0.436%
8	6.769	638	643	646	rVV	198091	285161	1.46%	0.434%
9	6.805	646	649	654	rVV	172274	253837	1.30%	0.386%
10	6.875	654	661	667	rVV3	281038	584885	2.99%	0.889%
11	6.934	667	671	678	rVB	116368	185986	0.95%	0.283%
12	7.099	694	699	703	rVV	115993	199596	1.02%	0.304%
13	7.140	703	706	710	rVV	134440	208352	1.07%	0.317%
14	7.334	733	739	748	rVV2	963860	1789438	9.16%	2.721%
15	7.622	778	788	793	rVV6	70651	203172	1.04%	0.309%
16	7.699	793	801	806	rVV2	484043	973668	4.98%	1.481%
17	7.764	806	812	817	rVV3	273845	476504	2.44%	0.725%
18	7.816	817	821	829	rVB3	112286	200038	1.02%	0.304%
19	7.987	847	850	857	rVB4	106469	184899	0.95%	0.281%
20	8.122	867	873	878	rBV	460638	739637	3.79%	1.125%
21	8.228	887	891	897	rVB2	148991	293152	1.50%	0.446%
22	8.281	897	900	904	rBV	109139	139332	0.71%	0.212%
23	8.334	904	909	910	rBV3	195017	292611	1.50%	0.445%
24	8.458	925	930	933	rBV	198313	285996	1.46%	0.435%
25	8.493	933	936	945	rVB2	126466	222389	1.14%	0.338%
26	8.640	956	961	965	rBV	117827	171386	0.88%	0.261%
27	8.687	965	969	977	rVB2	162506	295972	1.52%	0.450%
28	8.793	977	987	993	rBV3	187253	466229	2.39%	0.709%
29	8.916	999	1008	1013	rVB	924792	1441228	7.38%	2.192%
30	9.034	1020	1028	1034	rVB7	123531	335421	1.72%	0.510%
31	9.469	1095	1102	1108	rVB5	56742	144363	0.74%	0.220%
32	9.558	1109	1117	1121	rBV6	99068	240940	1.23%	0.366%
33	9.616	1121	1127	1131	rBV5	102104	191115	0.98%	0.291%
34	9.752	1147	1150	1156	rVB2	117543	193219	0.99%	0.294%
35	9.822	1157	1162	1167	rBV3	92504	151277	0.77%	0.230%
36	10.163	1214	1220	1225	rBV2	143752	243413	1.25%	0.370%

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Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

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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
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37	10.458	1259	1270	1277	rBV4	1346386	3016085	15.44%	4.587%
38	10.587	1286	1292	1296	rBV3	373166	637075	3.26%	0.969%
39	10.628	1296	1299	1305	rVB2	107303	139349	0.71%	0.212%
40	10.746	1314	1319	1329	rVB6	59987	134410	0.69%	0.204%

41	10.840	1329	1335	1348	rBV	477197	863178	4.42%	1.313%
42	10.969	1348	1357	1363	rVV2	125381	255203	1.31%	0.388%
43	11.375	1417	1426	1436	rVB8	109454	310253	1.59%	0.472%
44	11.487	1440	1445	1450	rVV	184507	334046	1.71%	0.508%
45	11.552	1451	1456	1464	rVB	343235	608520	3.12%	0.925%

46	11.728	1477	1486	1495	rBV2	113643	260519	1.33%	0.396%
47	11.887	1504	1513	1517	rBV2	476208	766665	3.92%	1.166%
48	11.922	1517	1519	1524	rVB	172447	214526	1.10%	0.326%
49	12.128	1548	1554	1561	rVB2	372413	744811	3.81%	1.133%
50	12.299	1577	1583	1586	rBV5	122427	203639	1.04%	0.310%

51	12.852	1672	1677	1680	rBV2	108688	171563	0.88%	0.261%
52	13.134	1720	1725	1732	rVB	399625	625958	3.20%	0.952%
53	13.357	1760	1763	1770	rVB4	93995	141083	0.72%	0.215%
54	13.487	1776	1785	1790	rBV3	136123	348986	1.79%	0.531%
55	13.622	1804	1808	1812	rBV	161444	287869	1.47%	0.438%

56	13.793	1829	1837	1842	rVB2	149376	271919	1.39%	0.414%
57	13.863	1842	1849	1853	rBV3	97148	211442	1.08%	0.322%
58	13.946	1859	1863	1868	rBV	168920	228621	1.17%	0.348%
59	14.004	1869	1873	1876	rVV	81659	132480	0.68%	0.201%
60	14.228	1905	1911	1916	rBV	627383	888690	4.55%	1.352%

61	14.316	1919	1926	1932	rVV	975814	1575044	8.06%	2.395%
62	14.398	1933	1940	1945	rVV8	115220	246518	1.26%	0.375%
63	14.457	1945	1950	1957	rVB	511340	762451	3.90%	1.160%
64	14.540	1958	1964	1967	rBV4	90795	184368	0.94%	0.280%
65	14.722	1991	1995	2000	rVB3	98330	158488	0.81%	0.241%

66	14.793	2004	2007	2014	rVB2	88261	130947	0.67%	0.199%
67	14.857	2014	2018	2022	rBV5	131048	219635	1.12%	0.334%
68	14.898	2022	2025	2027	rVV2	93526	129480	0.66%	0.197%
69	14.951	2029	2034	2040	rVV	1084498	1761024	9.02%	2.678%
70	15.010	2042	2044	2049	rVV3	155586	184741	0.95%	0.281%

71	15.081	2051	2056	2065	rVB	529241	869585	4.45%	1.322%
72	15.187	2070	2074	2080	rVB	391773	553580	2.83%	0.842%
73	15.322	2093	2097	2101	rVB	164014	217617	1.11%	0.331%
74	15.428	2109	2115	2120	rVB2	192846	296518	1.52%	0.451%
75	15.610	2140	2146	2150	rBV	199173	296125	1.52%	0.450%

76	15.698	2156	2161	2166	rVB2	193796	304814	1.56%	0.464%
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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

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77	15.981	2204	2209	2213	rVV	239209	359979	1.84%	0.547%
78	16.069	2219	2224	2227	rBV	413927	675323	3.46%	1.027%
79	16.416	2279	2283	2287	rBV2	146825	199962	1.02%	0.304%
80	16.763	2335	2342	2355	rBV	12459787	19533117	100.00%	29.706%
81	16.881	2358	2362	2367	rVV	397786	525355	2.69%	0.799%
82	16.934	2367	2371	2381	rVB2	219631	407949	2.09%	0.620%
83	17.110	2396	2401	2406	rBV	1425750	1904549	9.75%	2.896%
84	17.216	2414	2419	2423	rVB	202095	318392	1.63%	0.484%
85	17.263	2424	2427	2432	rBV3	112607	153514	0.79%	0.233%
86	17.416	2449	2453	2461	rVB5	173834	324449	1.66%	0.493%
87	17.645	2486	2492	2496	rBV	314366	460772	2.36%	0.701%
88	17.710	2499	2503	2508	rVV3	78073	138826	0.71%	0.211%
89	17.816	2517	2521	2524	rVB	169387	195832	1.00%	0.298%
90	18.363	2610	2614	2619	rBV	284482	363289	1.86%	0.552%
91	19.022	2722	2726	2728	rBV	241572	307231	1.57%	0.467%
92	19.592	2819	2823	2826	rBV	256209	325967	1.67%	0.496%
93	19.633	2826	2830	2835	rVB2	150782	191035	0.98%	0.291%
94	20.181	2919	2923	2927	rBV2	228138	328608	1.68%	0.500%
95	20.704	3009	3012	3016	rBV	130826	140634	0.72%	0.214%
96	21.228	3096	3101	3108	rVB	392334	500639	2.56%	0.761%
97	21.539	3149	3154	3163	rBV2	1356193	2306627	11.81%	3.508%
98	22.286	3277	3281	3288	rVB4	161125	259362	1.33%	0.394%
99	24.604	3671	3675	3686	rVB	127917	317812	1.63%	0.483%
100	24.839	3705	3715	3727	rVB	1029838	2534110	12.97%	3.854%

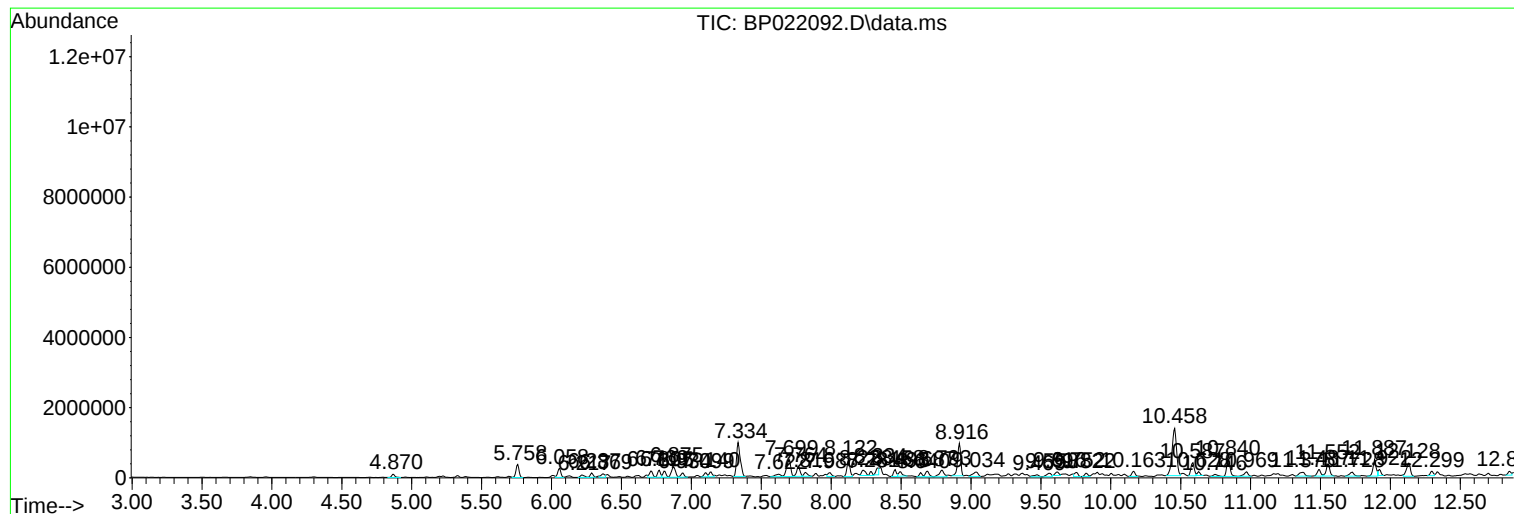
Sum of corrected areas: 65754426

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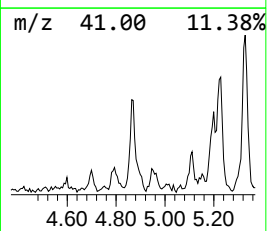
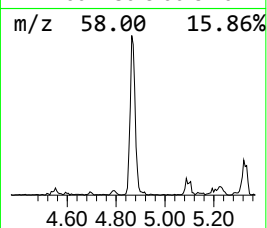
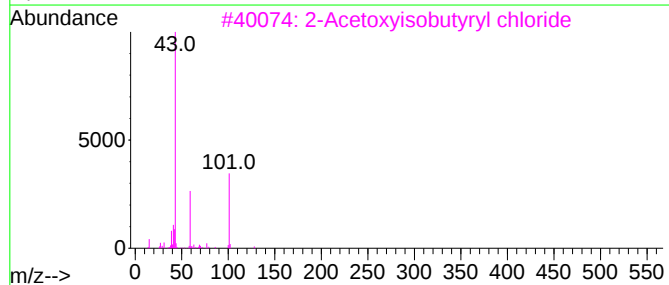
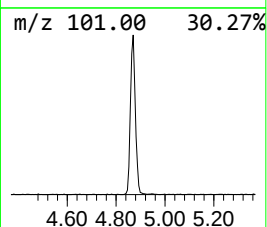
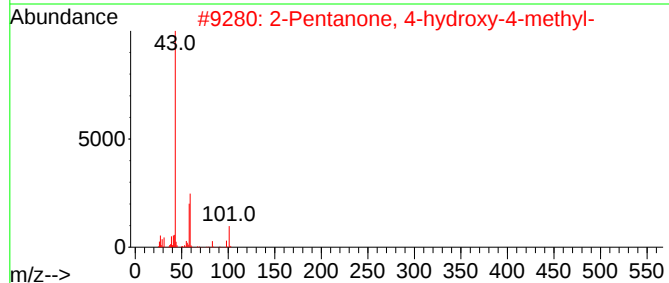
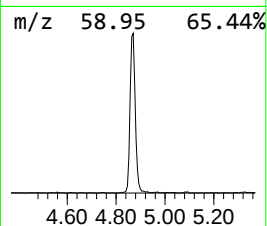
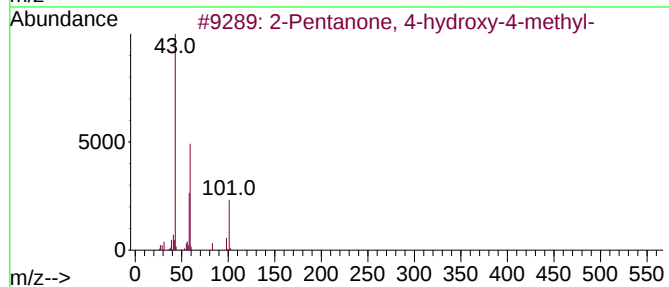
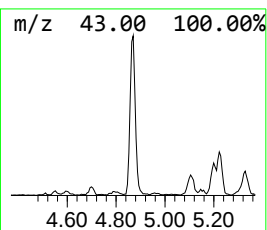
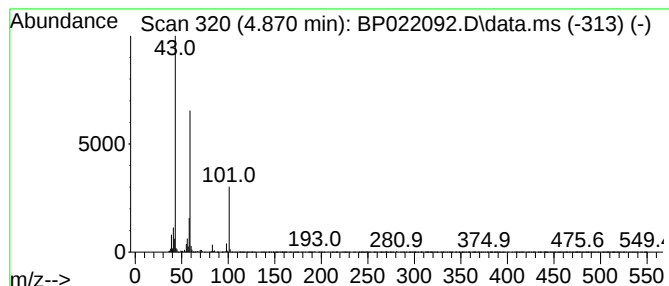
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.870	3.03 ng/ul	147328	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45		
3	2-Acetoxyisobutyryl chloride	164	C6H9ClO3	040635-66-3	38		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		
5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		



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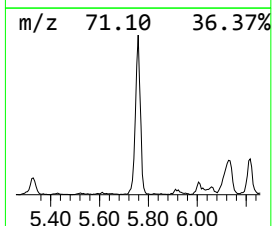
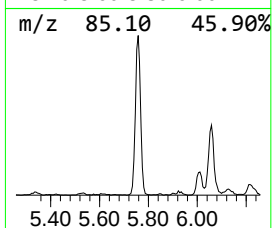
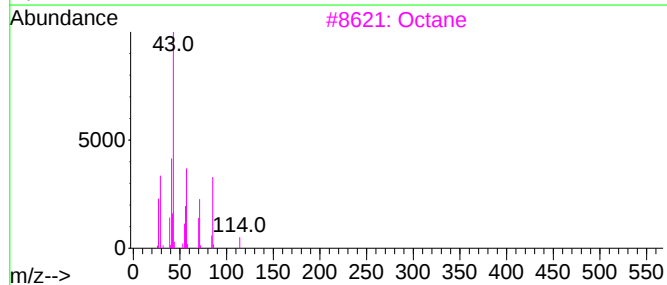
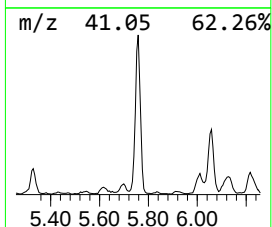
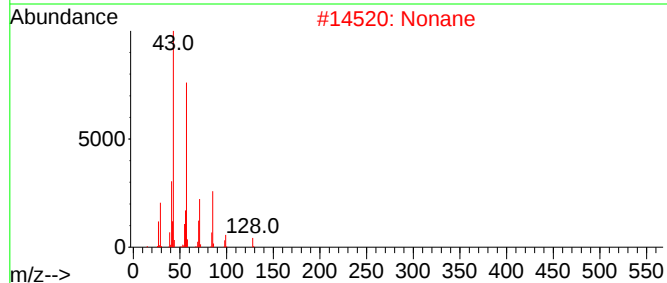
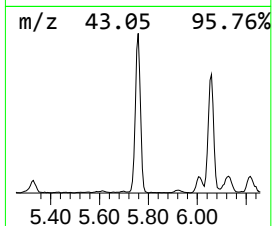
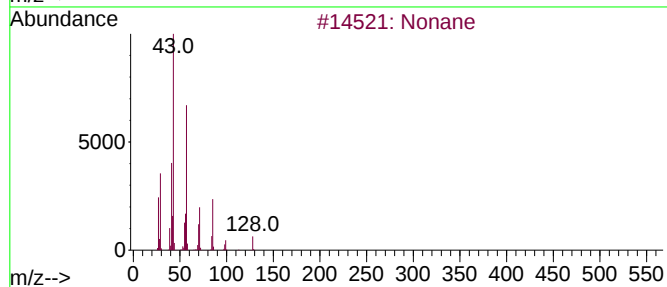
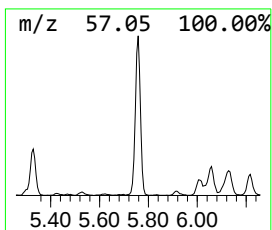
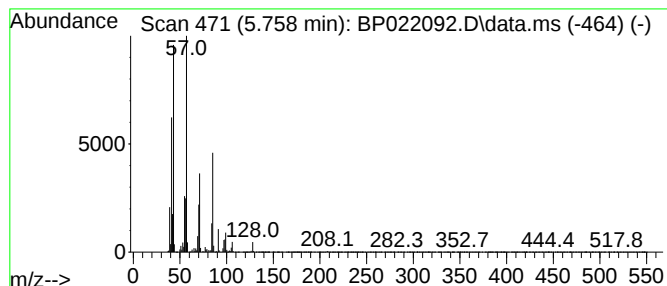
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.758	11.47 ng/ul	558495	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	94
2			Nonane	128	C9H20	000111-84-2	94
3			Octane	114	C8H18	000111-65-9	72
4			Octane	114	C8H18	000111-65-9	64
5			Decane	142	C10H22	000124-18-5	64



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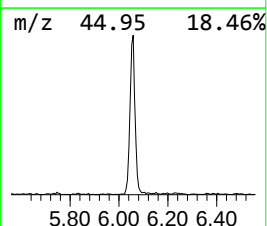
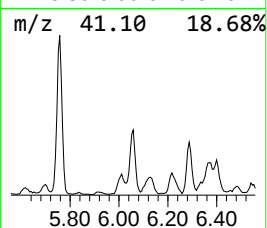
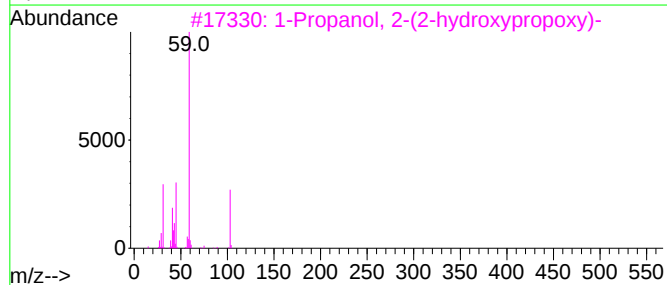
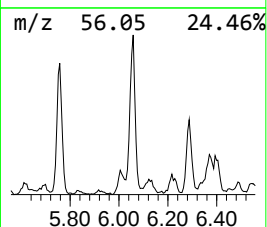
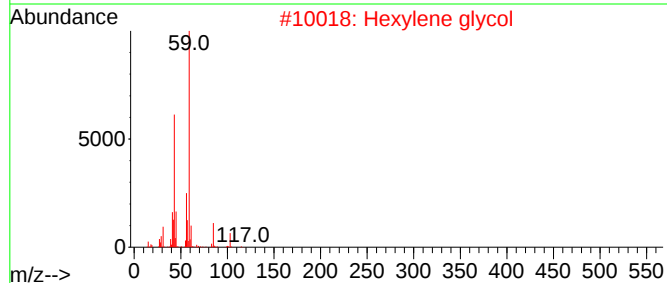
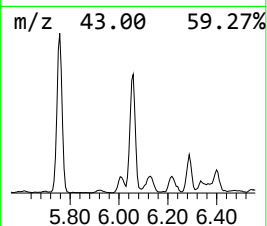
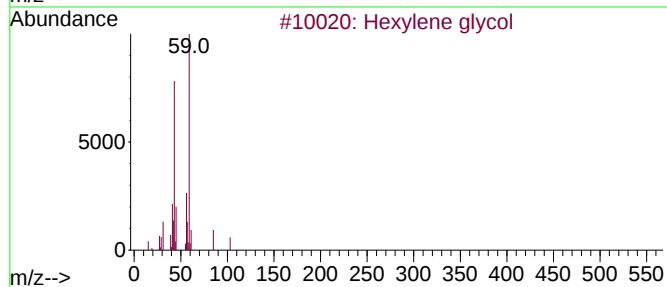
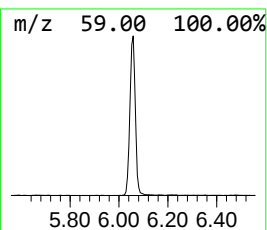
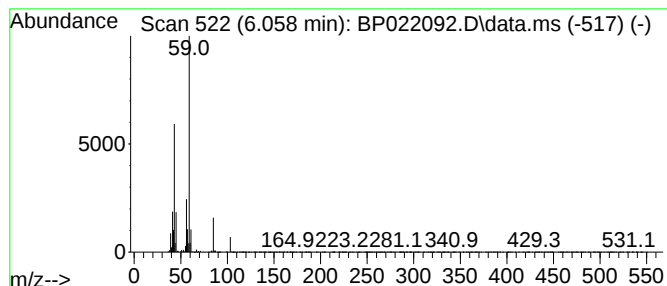
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Hexylene glycol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.058	8.10 ng/ul	394480	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexylene glycol	118	C6H14O2	000107-41-5	90
2			Hexylene glycol	118	C6H14O2	000107-41-5	83
3			1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	59
4			1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	59
5			Hexylene glycol	118	C6H14O2	000107-41-5	53



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Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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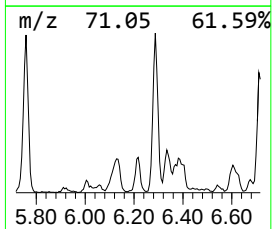
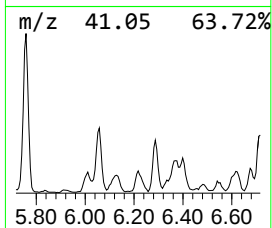
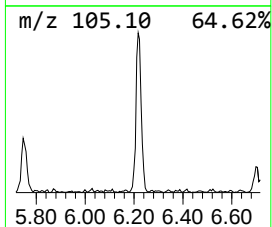
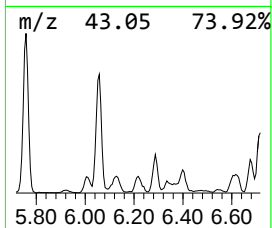
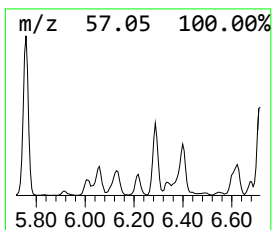
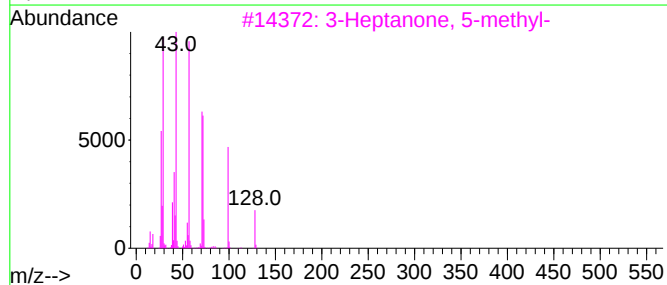
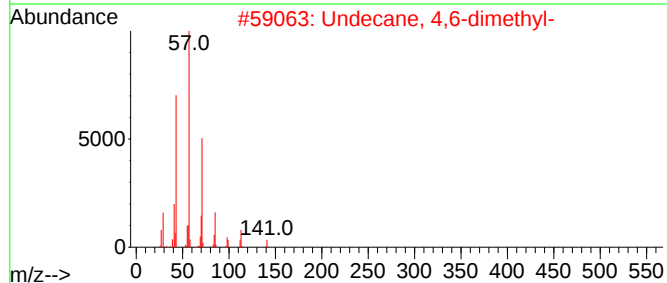
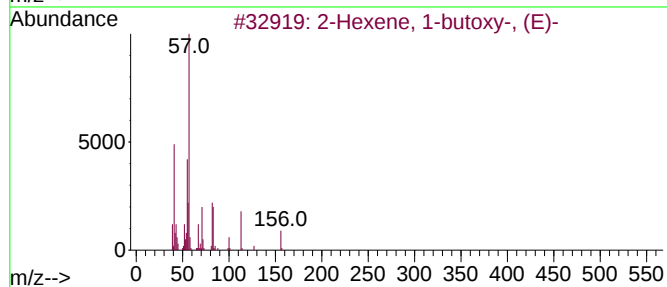
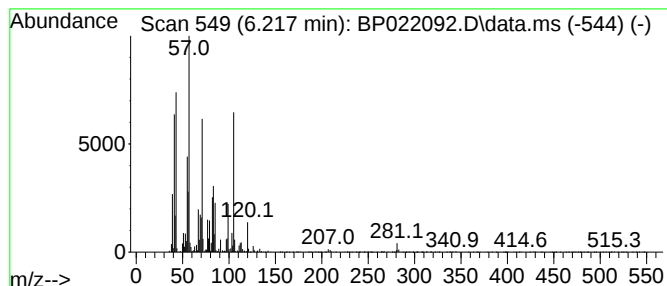
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-01 Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.217	2.80 ng/ul	136498	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexene, 1-butoxy-, (E)-		156	C10H20O	054340-67-9	38	
2	Undecane, 4,6-dimethyl-		184	C13H28	017312-82-2	35	
3	3-Heptanone, 5-methyl-		128	C8H16O	000541-85-5	35	
4	Tridecane		184	C13H28	000629-50-5	35	
5	Oxalic acid, hexyl octadecyl ester		426	C26H50O4	1000309-25-2	35	



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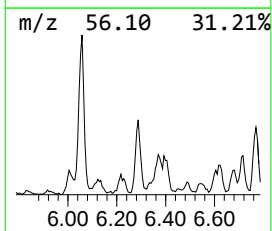
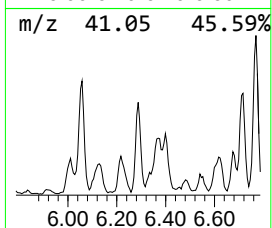
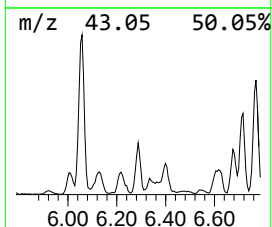
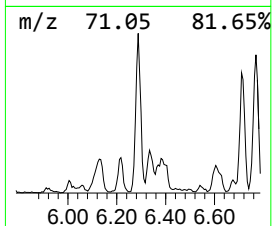
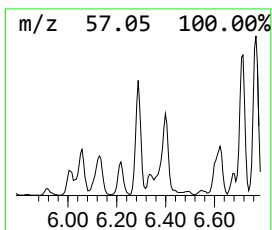
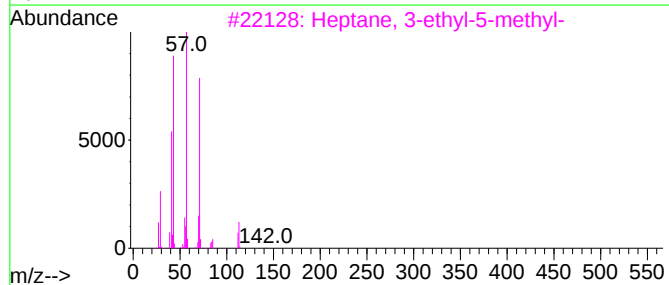
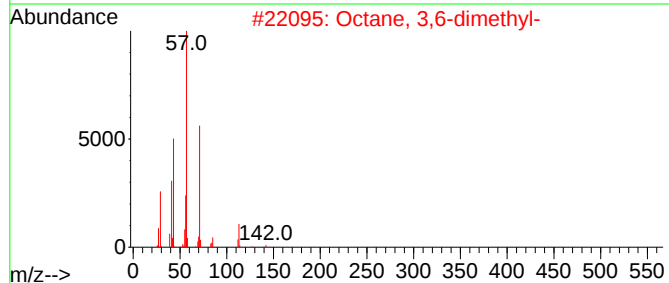
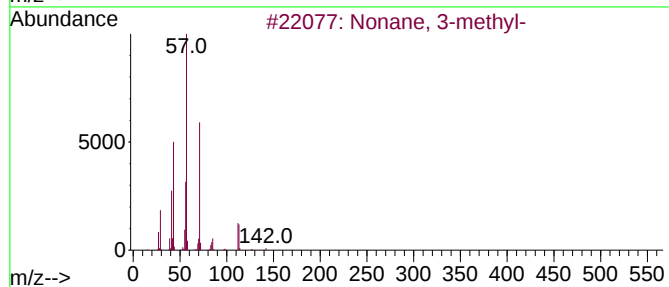
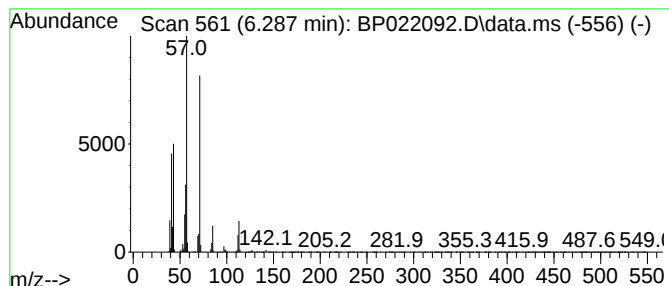
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.287	3.67 ng/ul	178818	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3-methyl-	142	C10H22	005911-04-6	91
2			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	90
3			Heptane, 3-ethyl-5-methyl-	142	C10H22	052896-90-9	72
4			Octane, 1,1'-oxybis-	242	C16H34O	000629-82-3	72
5			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
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Acq On : 28 Sep 2024 05:10
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Sample : P3910-17 10X
Misc :
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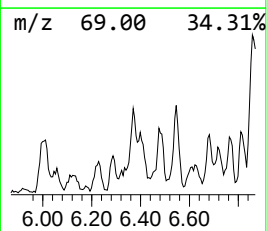
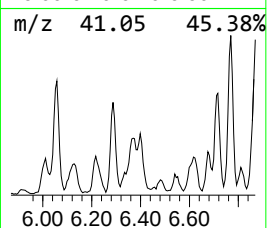
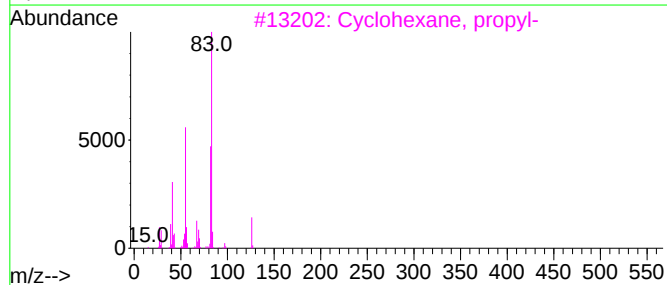
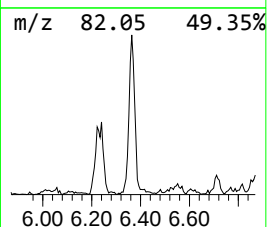
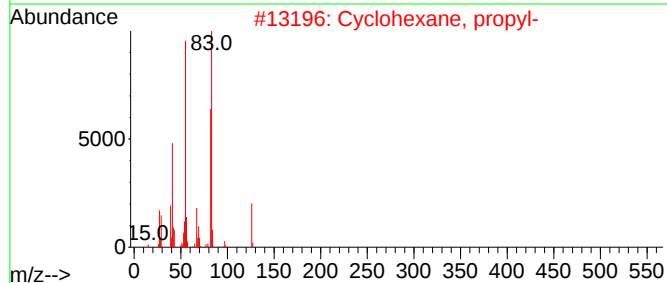
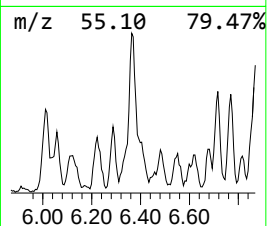
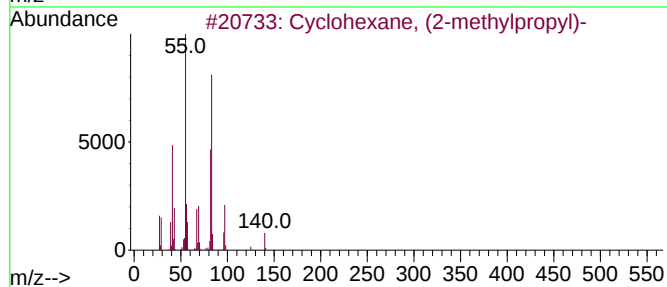
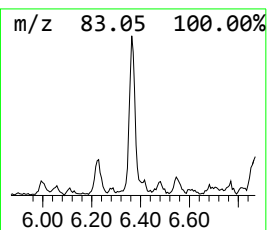
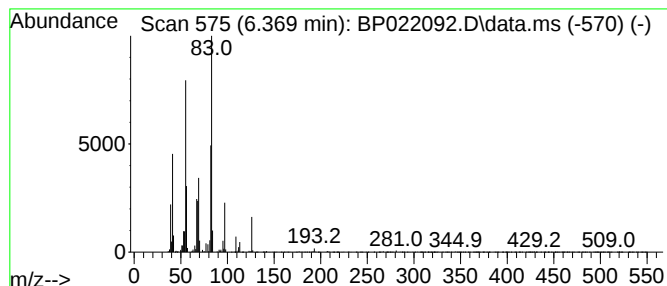
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic6.369 Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.369	4.07 ng/ul	197928	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	72
2			Cyclohexane, propyl-	126	C9H18	001678-92-8	64
3			Cyclohexane, propyl-	126	C9H18	001678-92-8	64
4			Cyclohexane, propyl-	126	C9H18	001678-92-8	64
5			Cyclohexane, (3-methylpentyl)-	168	C12H24	061142-38-9	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
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Acq On : 28 Sep 2024 05:10
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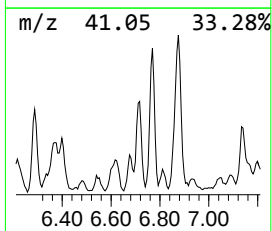
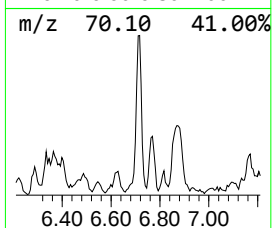
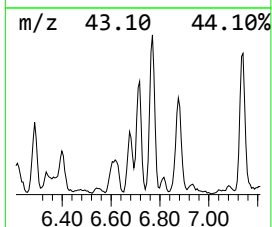
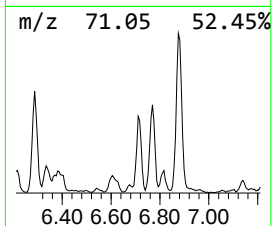
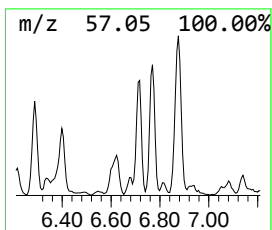
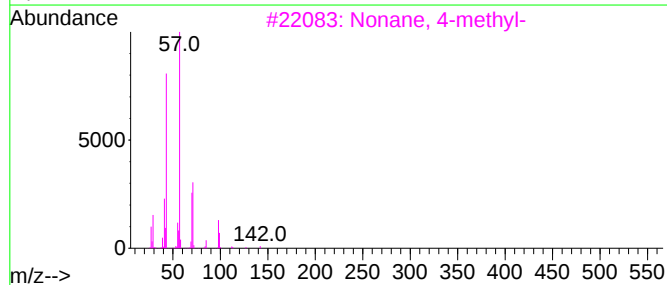
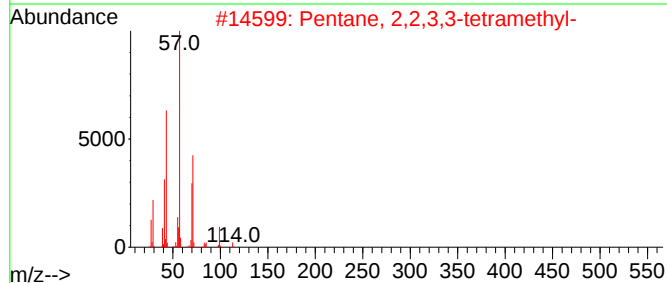
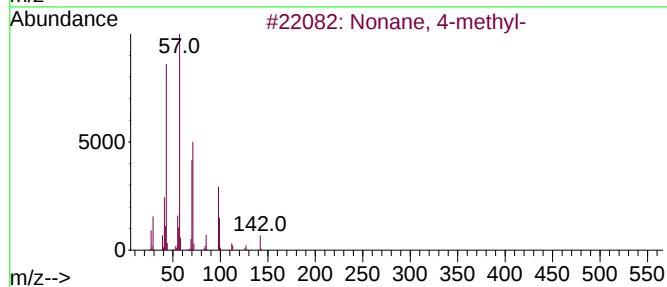
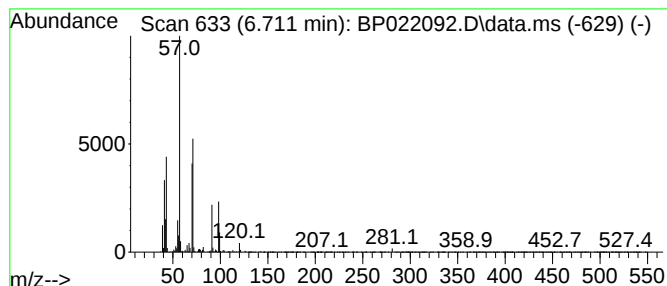
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.711	5.89 ng/ul	286505	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 4-methyl-	142	C10H22	017301-94-9	56
2			Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	53
3			Nonane, 4-methyl-	142	C10H22	017301-94-9	47
4			Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	45
5			Nonane, 4-methyl-	142	C10H22	017301-94-9	45



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Misc :
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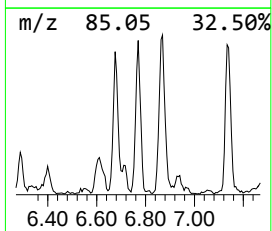
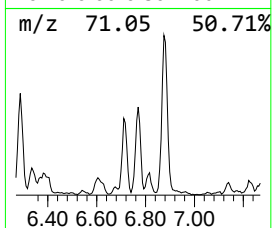
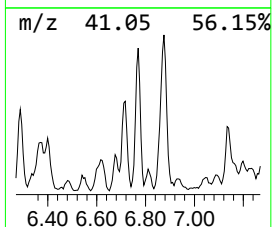
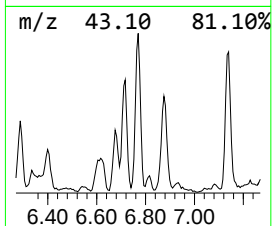
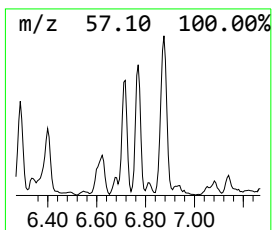
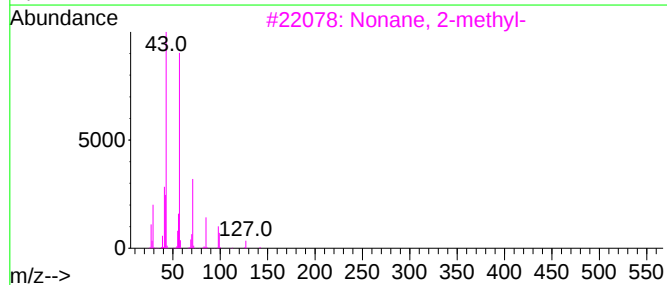
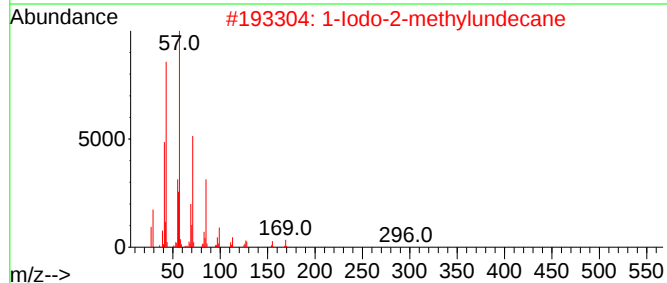
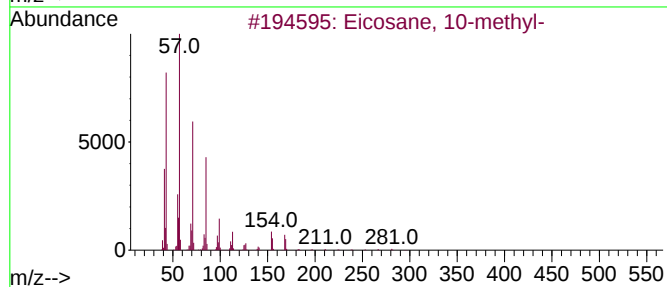
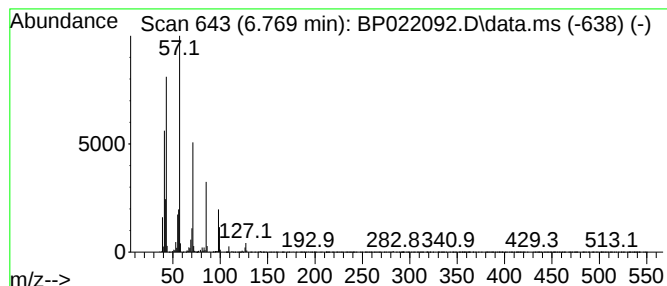
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.769	5.86 ng/ul	285161	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 10-methyl-	296	C21H44	054833-23-7	64		
2	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	64		
3	Nonane, 2-methyl-	142	C10H22	000871-83-0	58		
4	1-Iodoundecane	282	C11H23I	004282-44-4	55		
5	Tridecane	184	C13H28	000629-50-5	53		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
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ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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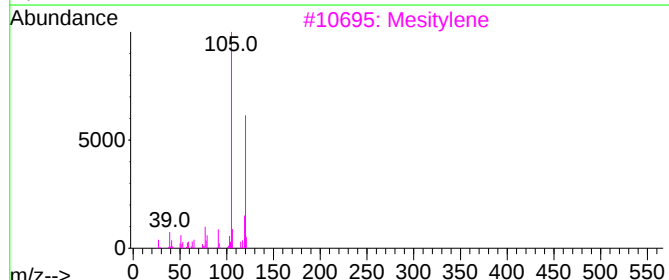
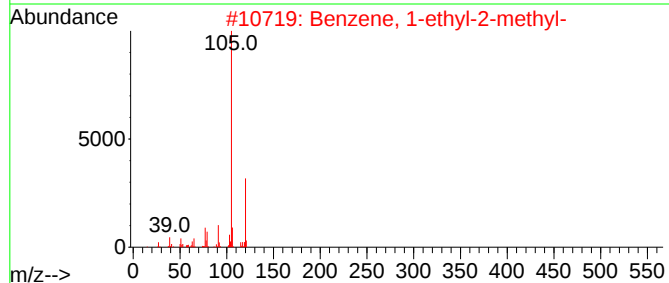
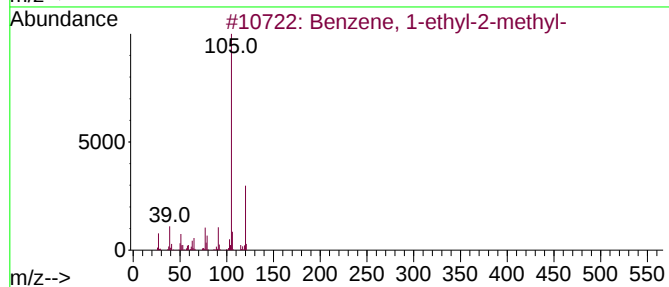
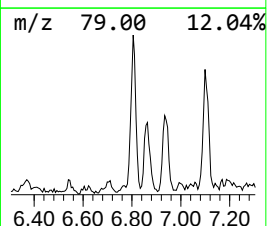
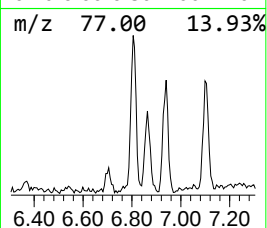
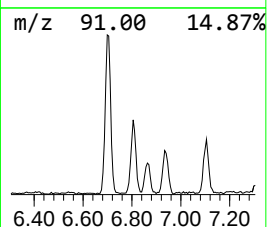
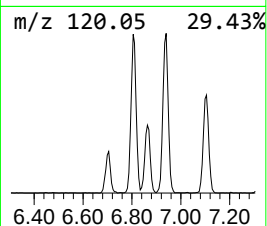
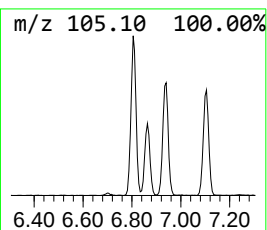
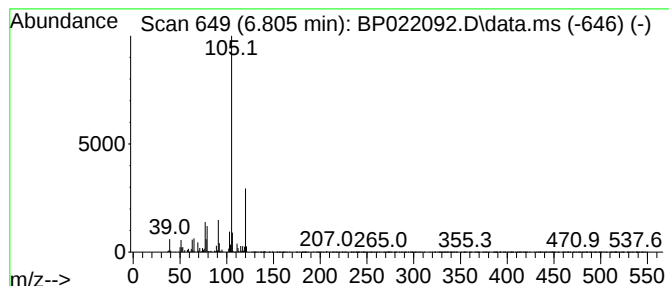
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.805	5.21 ng/ul	253837	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	94
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
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ALS Vial : 21 Sample Multiplier: 1

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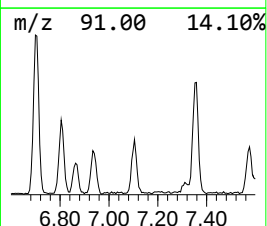
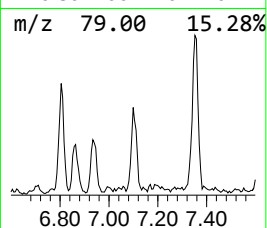
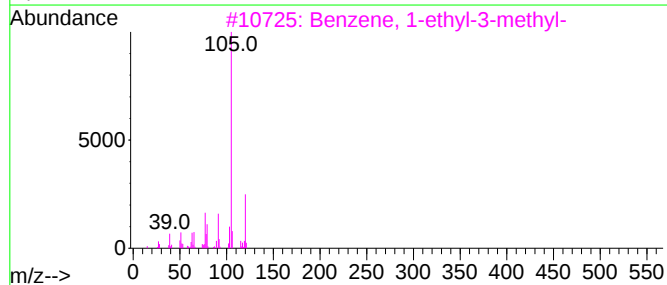
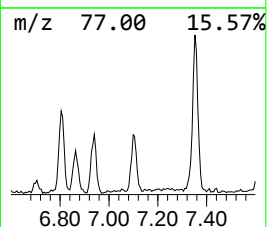
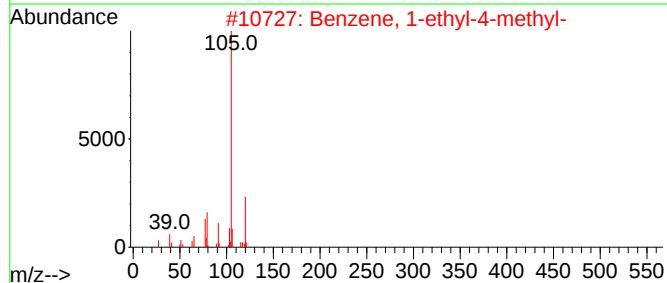
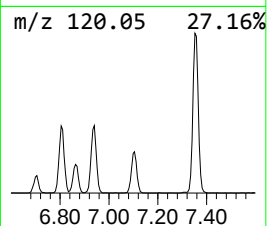
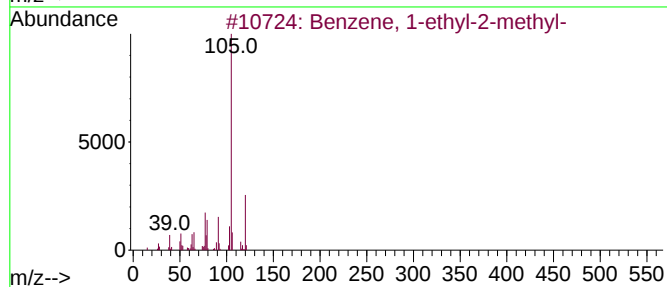
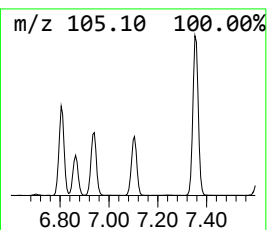
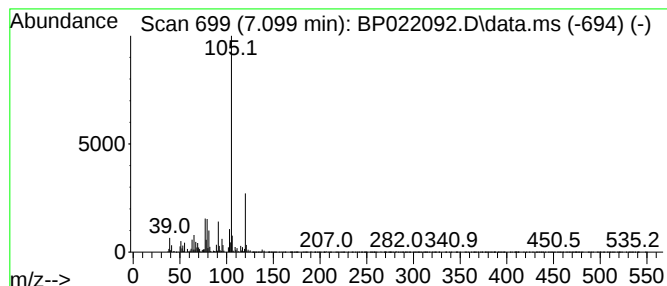
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzene, 1-ethyl-4-methyl- Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.099	4.10 ng/ul	199596	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	94
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	81
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	81
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC63

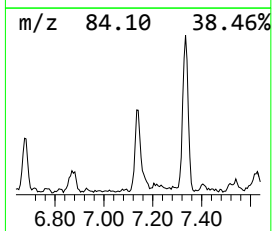
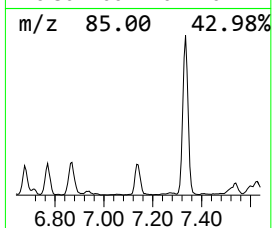
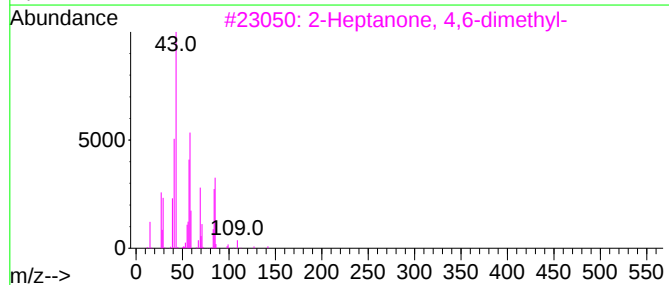
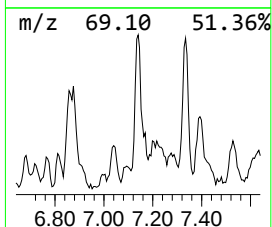
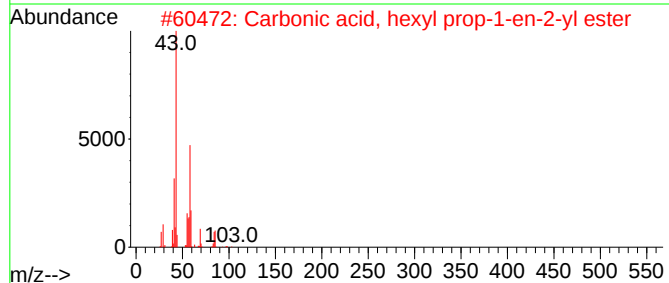
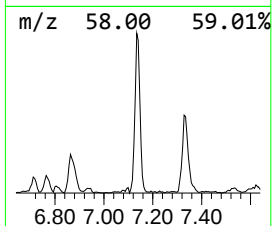
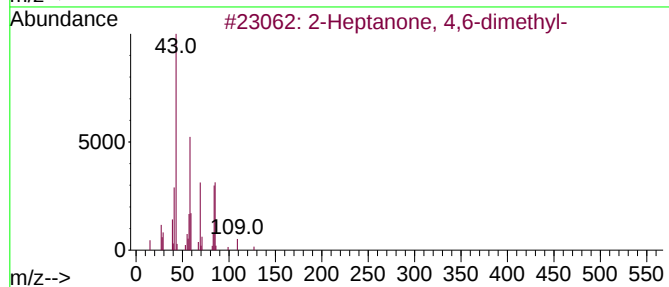
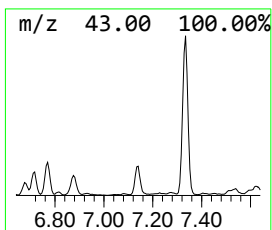
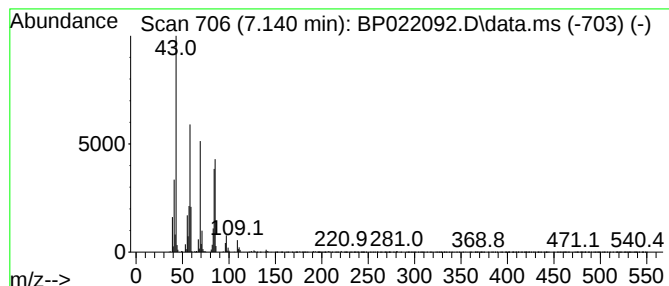
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 2-Heptanone, 4,6-dimethyl- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.140	4.28 ng/ul	208352	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	86
2			Carbonic acid, hexyl prop-1-en-2-yl ester	186	C10H18O3	1000382-54-2	72
3			2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	50
4			trans-2,3-Epoxydecane	156	C10H20O	054125-39-2	46
5			Nonane, 5-methyl-	142	C10H22	015869-85-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
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Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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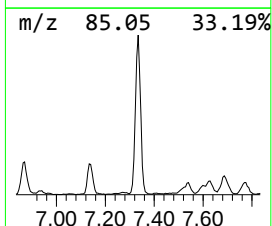
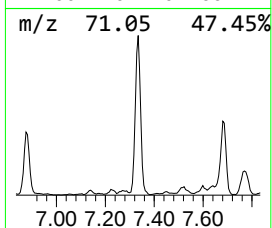
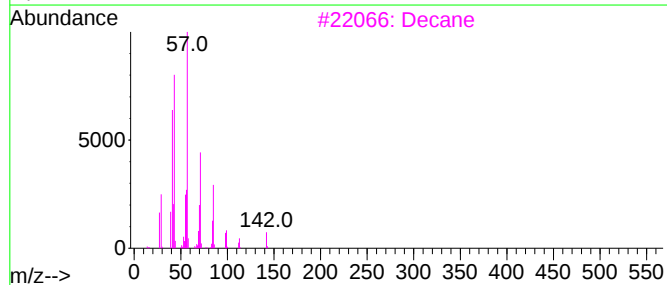
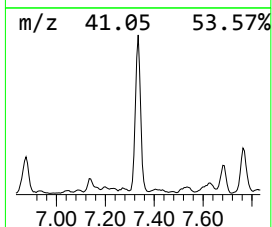
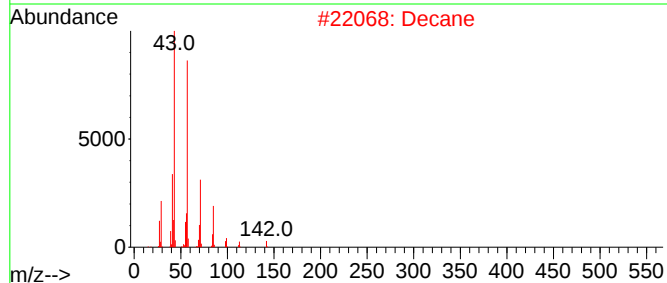
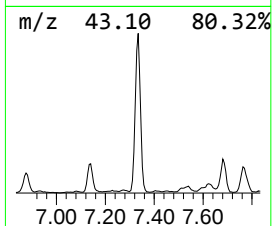
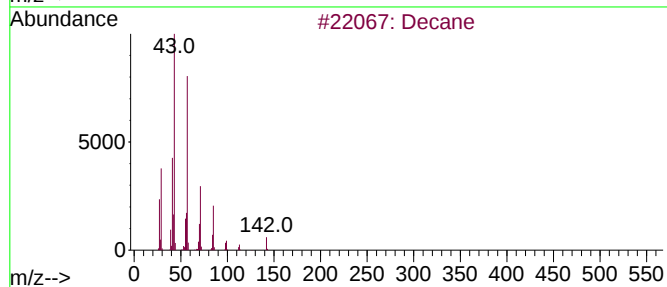
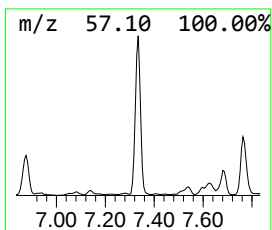
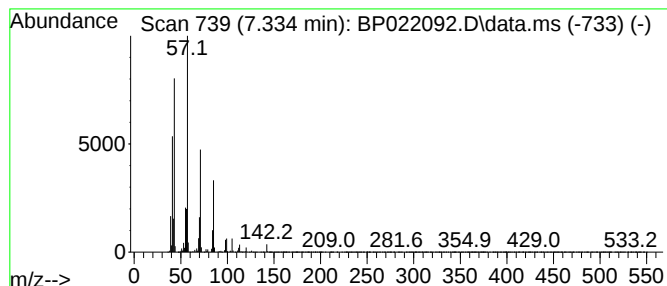
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.334	36.76 ng/ul	1789440	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	95
2			Decane	142	C10H22	000124-18-5	94
3			Decane	142	C10H22	000124-18-5	90
4			Tetradecane	198	C14H30	000629-59-4	78
5			Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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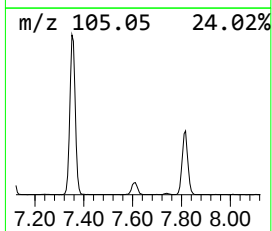
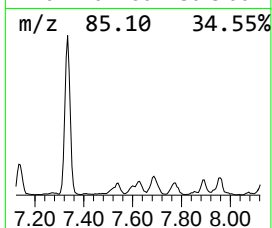
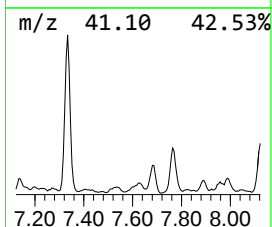
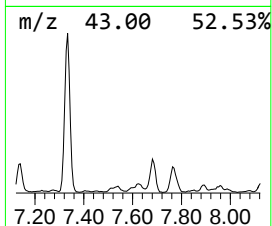
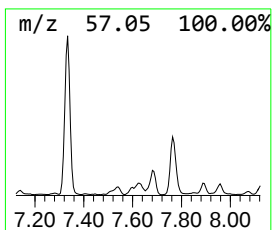
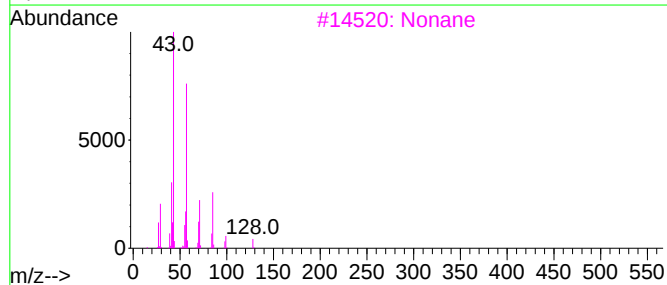
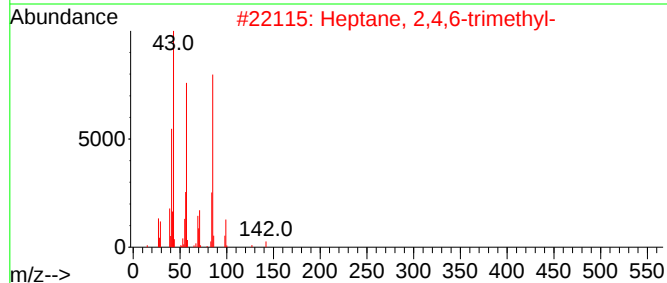
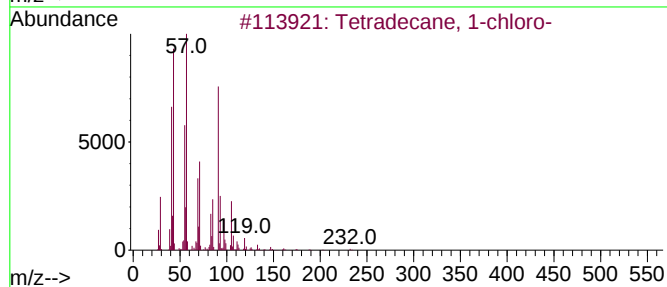
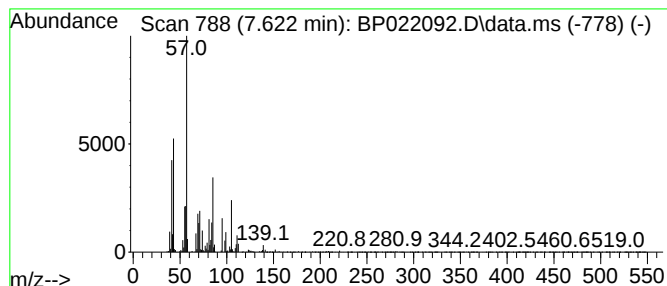
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Tetradecane, 1-chloro- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.622	4.17 ng/ul	203172	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane, 1-chloro-	232	C14H29Cl	002425-54-9	50
2			Heptane, 2,4,6-trimethyl-	142	C10H22	002613-61-8	46
3			Nonane	128	C9H20	000111-84-2	43
4			Nonane, 2,5-dimethyl-	156	C11H24	017302-27-1	43
5			Oxalic acid, hexyl pentadecyl ester	384	C23H44O4	1010309-24-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
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Misc :
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Instrument :
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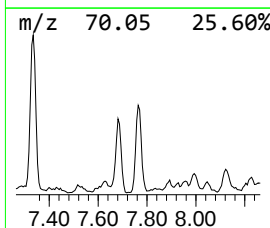
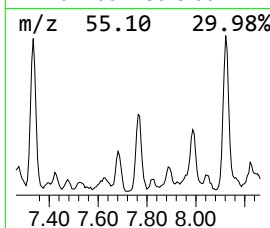
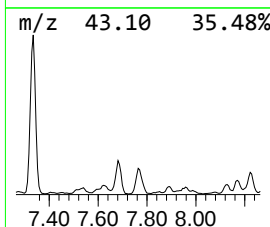
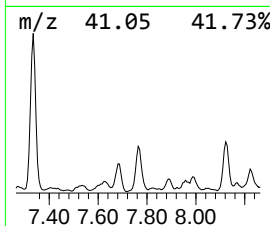
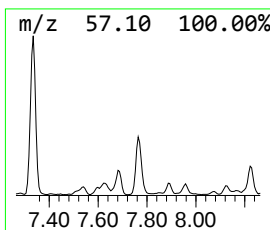
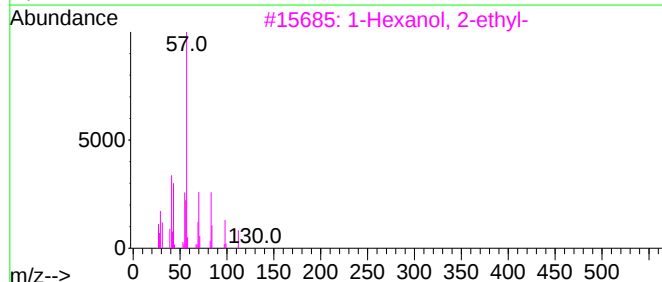
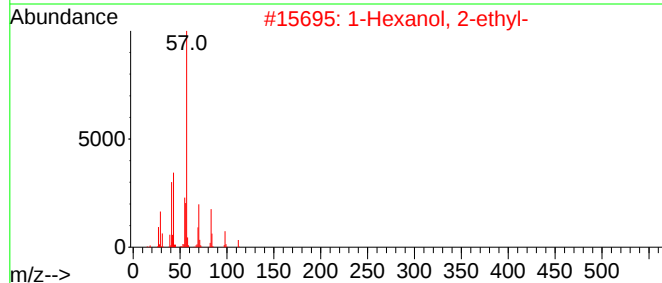
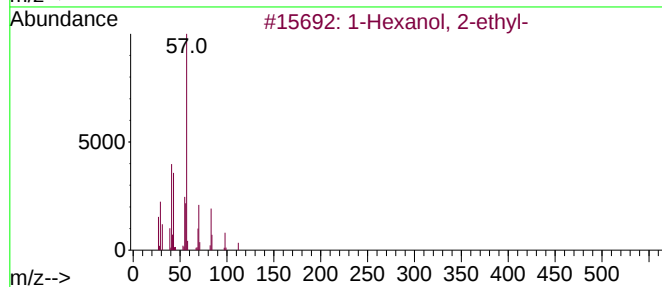
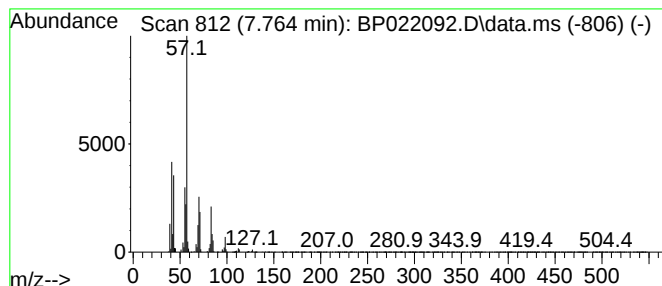
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 1-Hexanol, 2-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.764	9.79 ng/ul	476504	1,4-Dichlorobenzene-d4	7.699

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
2		1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72
3		1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64
4		1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	59
5		1-Decene, 2,4-dimethyl-	168	C12H24	055170-80-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
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Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

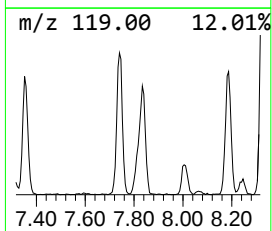
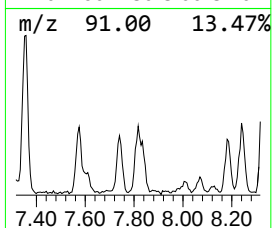
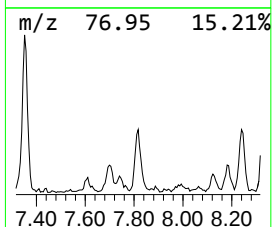
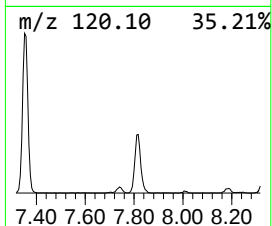
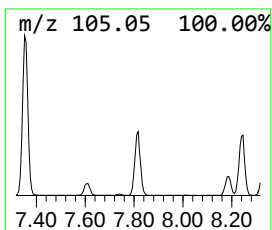
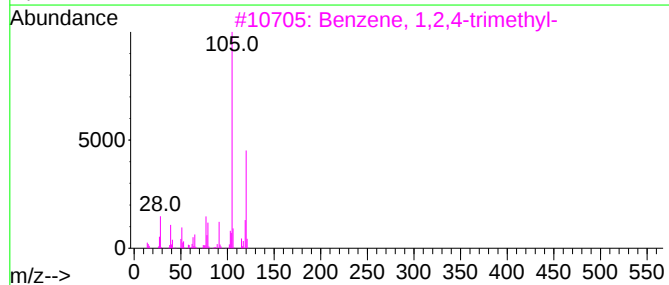
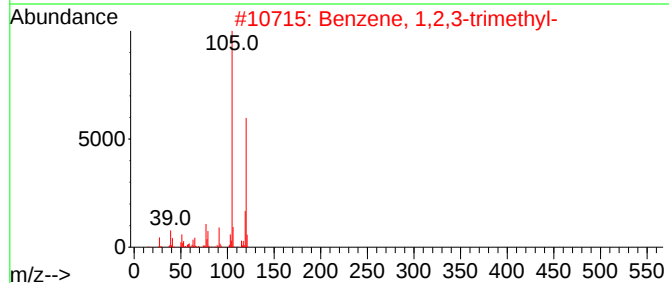
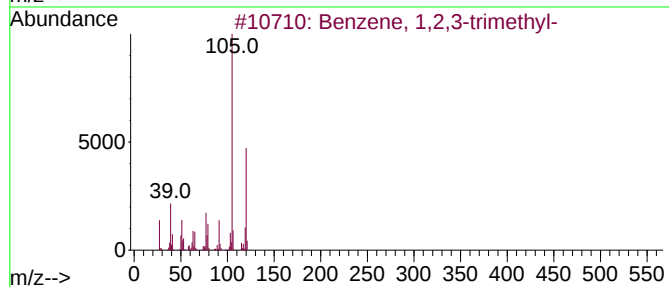
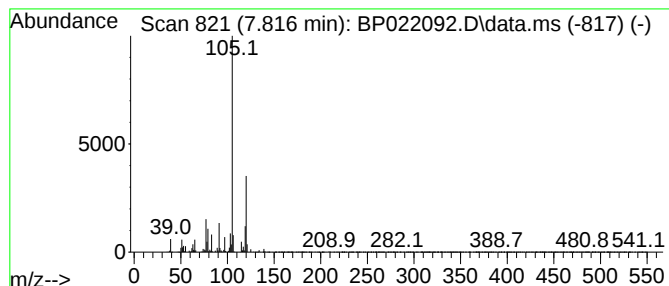
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Benzene, 1,2,3-trimethyl- Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD			R.T.
7.816	4.11 ng/ul	200038	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	94
2	Benzene, 1,2,3-trimethyl-		120	C9H12	000526-73-8	94
3	Benzene, 1,2,4-trimethyl-		120	C9H12	000095-63-6	93
4	Benzene, 1,2,3-trimethyl-		120	C9H12	000526-73-8	93
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 05:10
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ClientSampleId :
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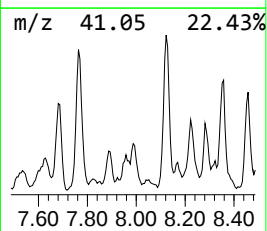
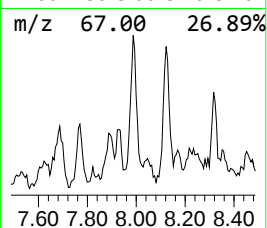
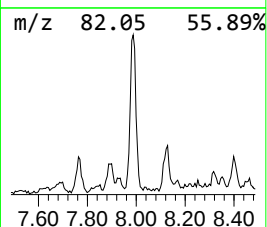
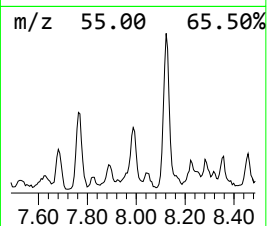
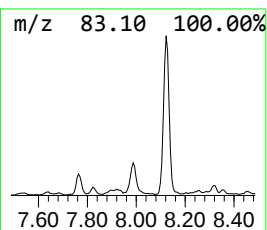
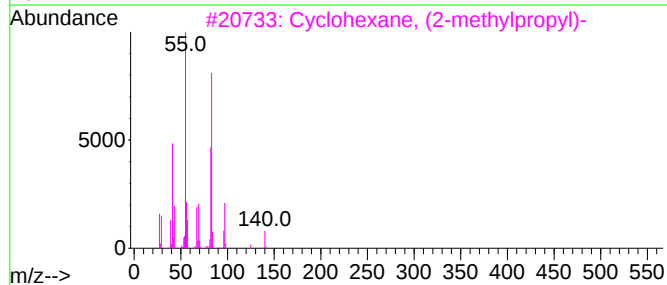
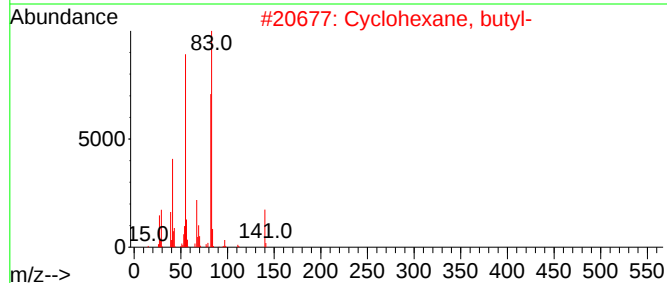
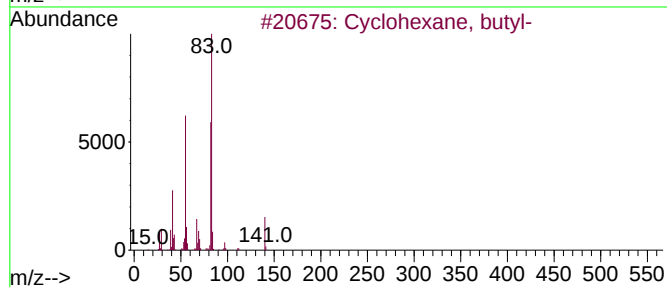
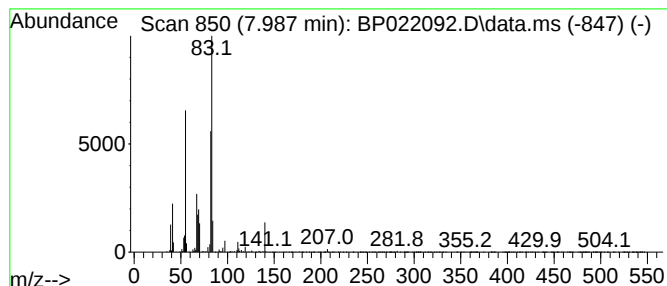
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Cyclic7.987 Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.987	3.80 ng/ul	184899	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, butyl-	140	C10H20	001678-93-9	83
2			Cyclohexane, butyl-	140	C10H20	001678-93-9	72
3			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	72
4			Cyclohexane, pentyl-	154	C11H22	004292-92-6	64
5			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC63

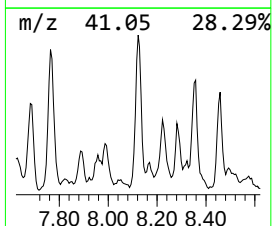
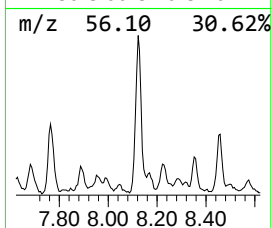
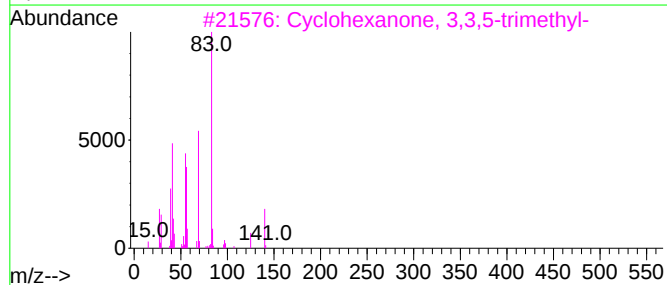
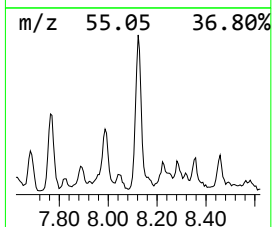
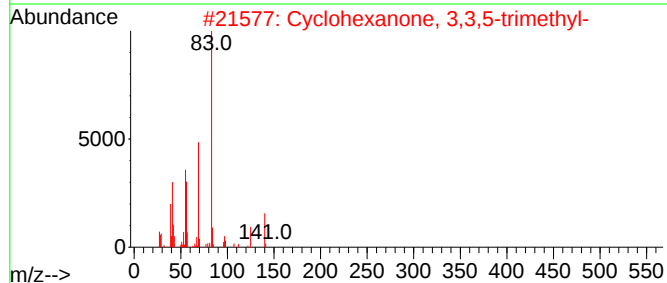
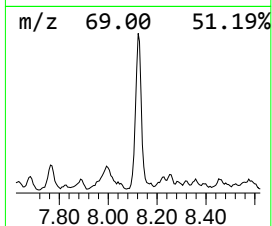
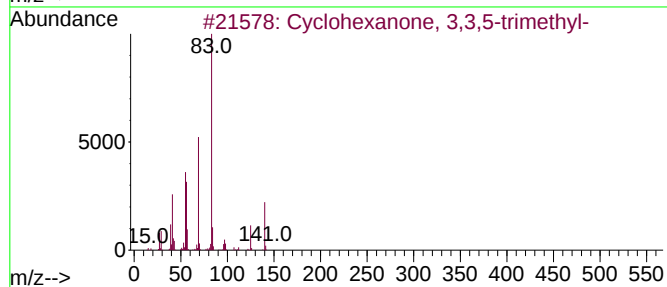
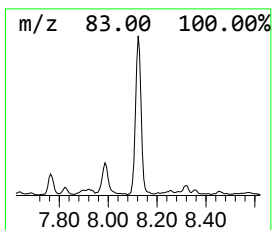
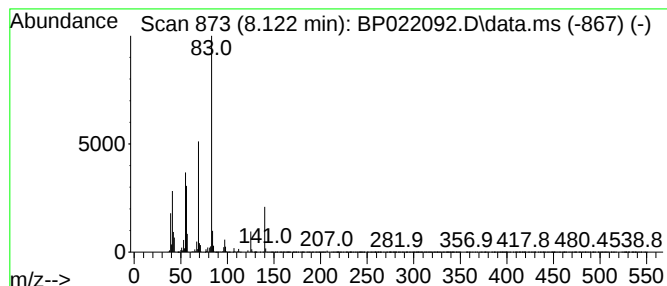
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Cyclohexanone, 3,3,5-trimet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.122	15.19 ng/ul	739637	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	96
2			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	95
3			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
4			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
5			3-Acetamidofuran	125	C6H7NO2	059445-85-1	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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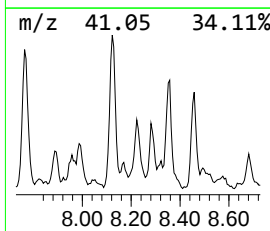
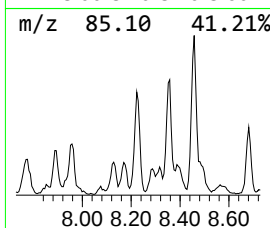
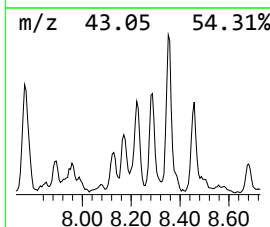
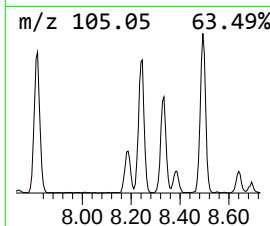
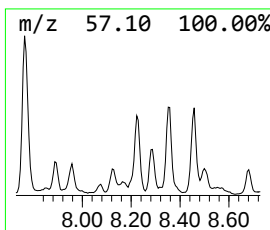
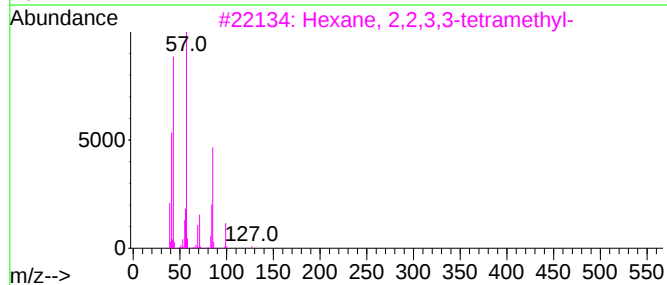
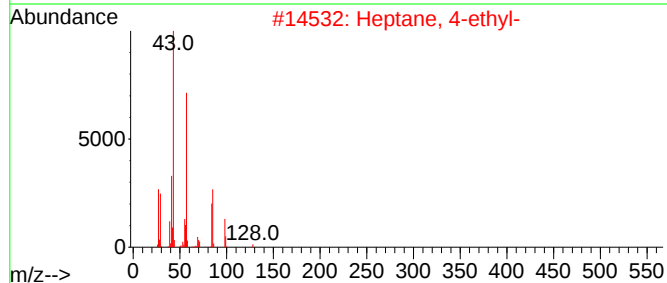
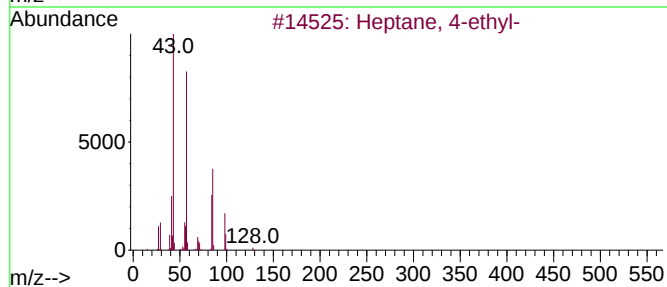
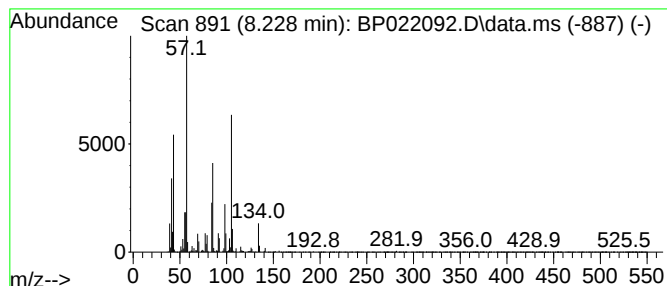
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.228	6.02 ng/ul	293152	1,4-Dichlorobenzene-d4	7.699

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 4-ethyl-	128	C9H20	002216-32-2	50
2		Heptane, 4-ethyl-	128	C9H20	002216-32-2	43
3		Hexane, 2,2,3,3-tetramethyl-	142	C10H22	013475-81-5	38
4		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	35
5		Pentane, 3-ethyl-2,3-dimethyl-	128	C9H20	016747-33-4	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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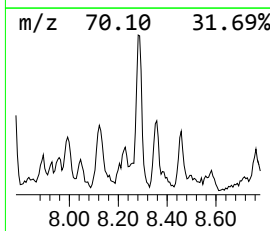
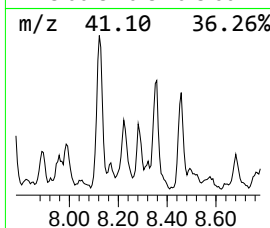
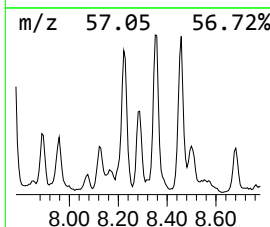
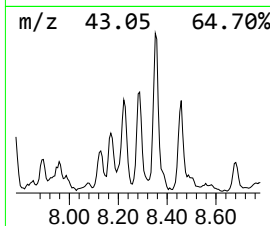
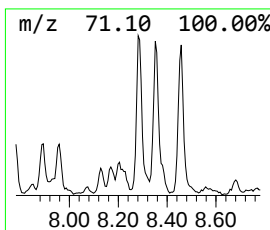
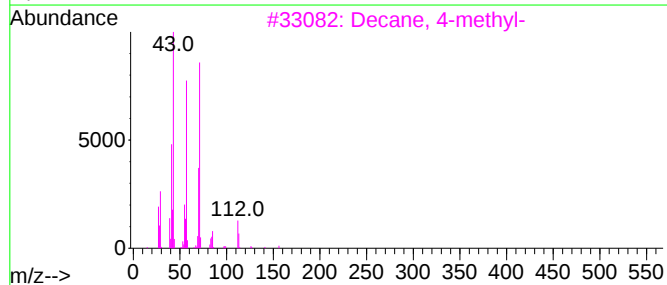
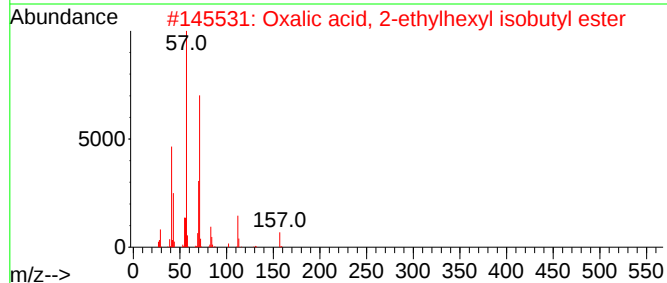
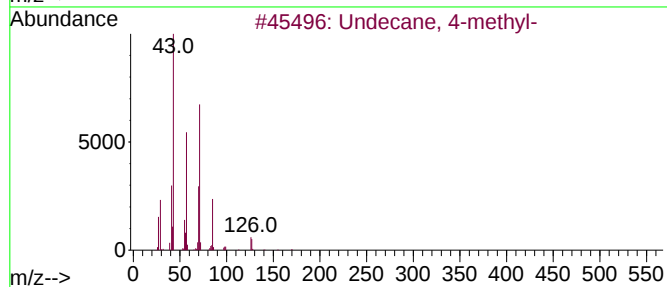
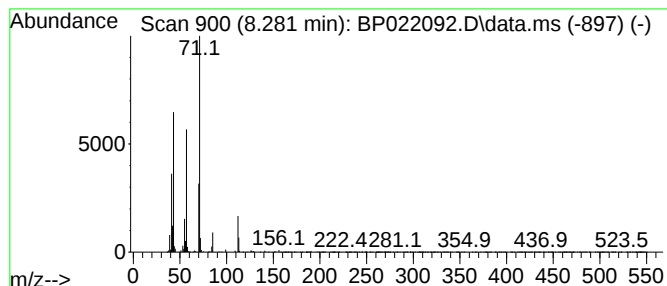
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.281	2.86 ng/ul	139332	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 4-methyl-	170	C12H26	002980-69-0	64
2			Oxalic acid, 2-ethylhexyl isobut...	258	C14H26O4	1010309-38-5	64
3			Decane, 4-methyl-	156	C11H24	002847-72-5	62
4			Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	59
5			Decane, 3,3,6-trimethyl-	184	C13H28	062338-14-1	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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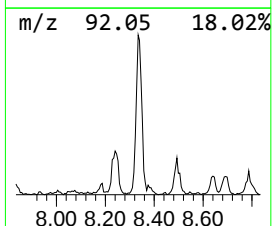
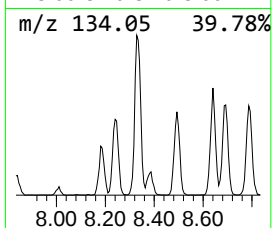
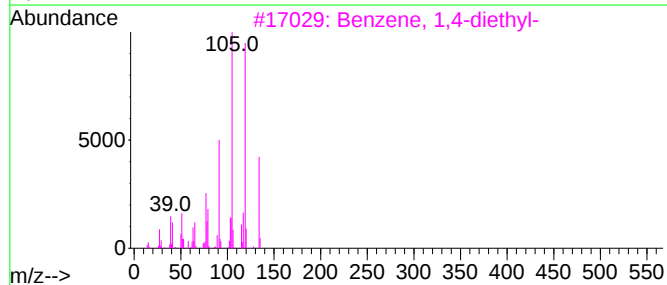
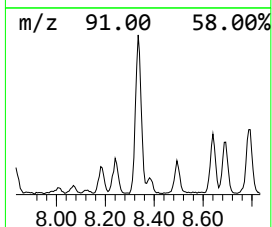
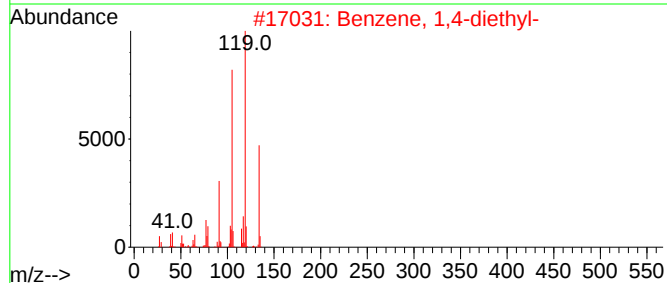
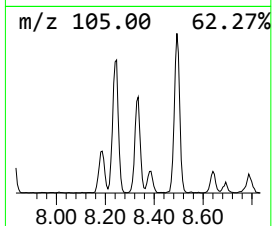
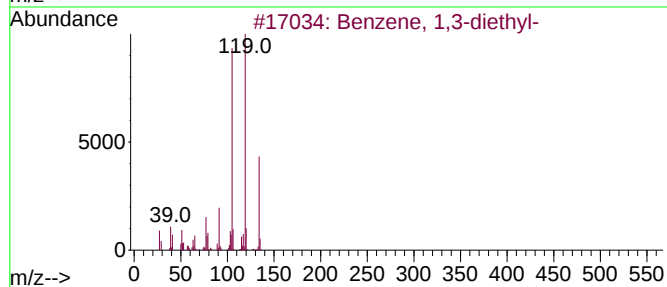
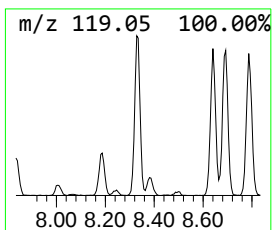
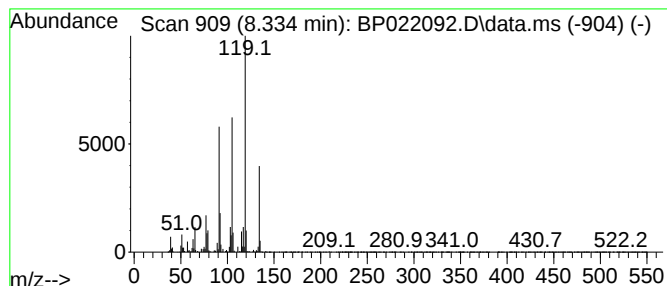
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Benzene, 1,3-diethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.334	6.01 ng/ul	292611	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
2			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
4			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	94
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
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Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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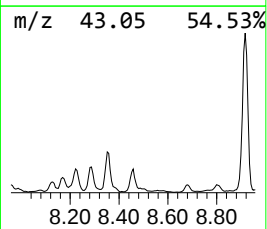
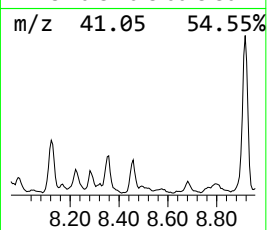
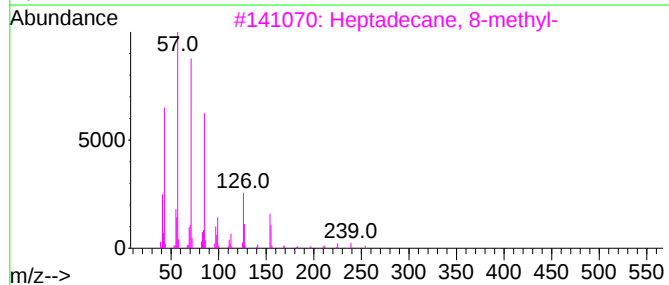
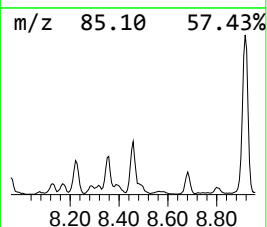
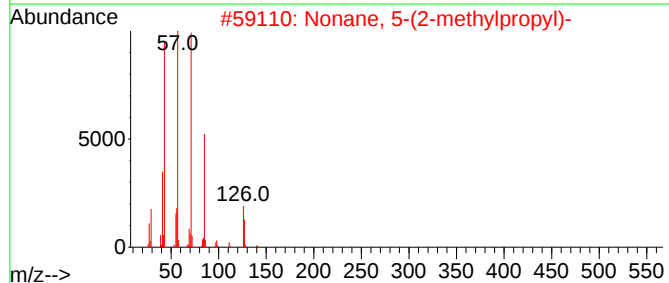
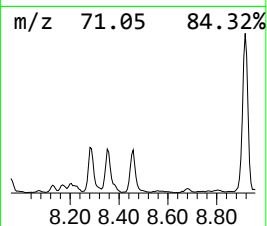
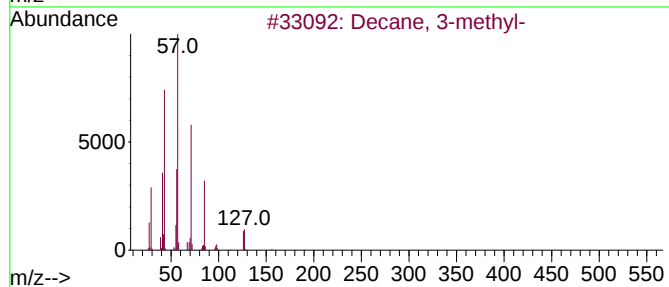
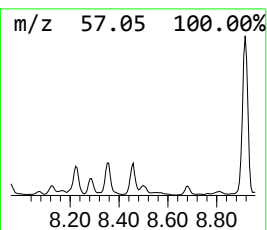
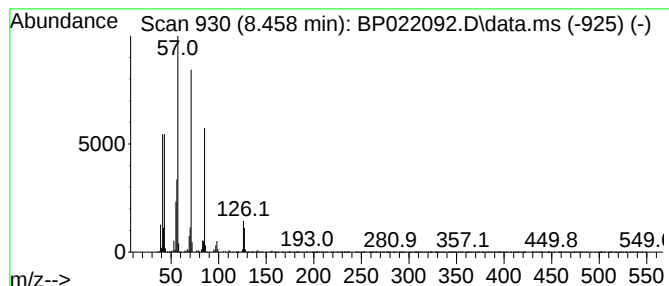
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.458	5.87 ng/ul	285996	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, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	78
3		Heptadecane, 8-methyl-	254	C18H38	013287-23-5	72
4		Decane, 3-methyl-	156	C11H24	013151-34-3	64
5		Nonane, 1-iodo-	254	C9H19I	004282-42-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
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Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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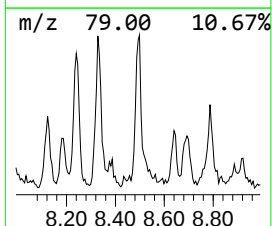
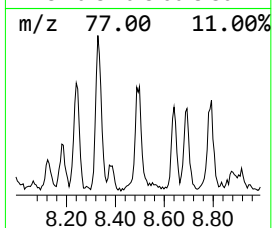
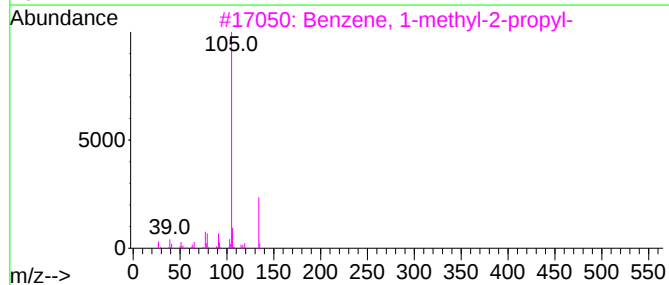
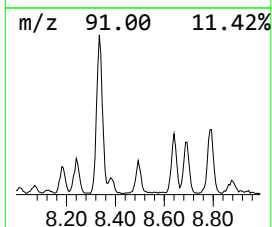
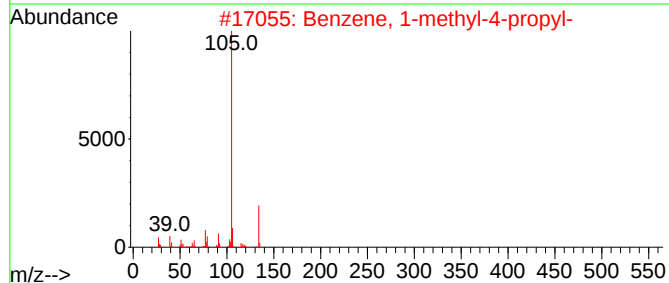
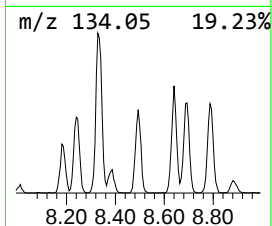
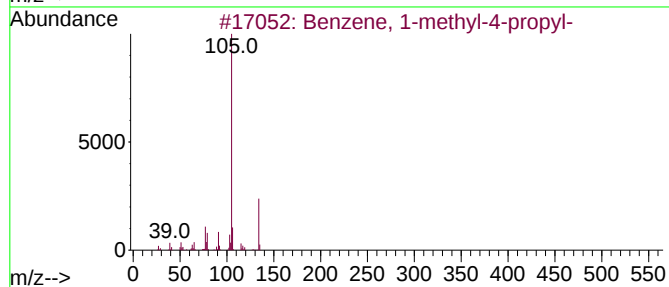
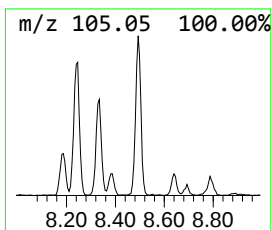
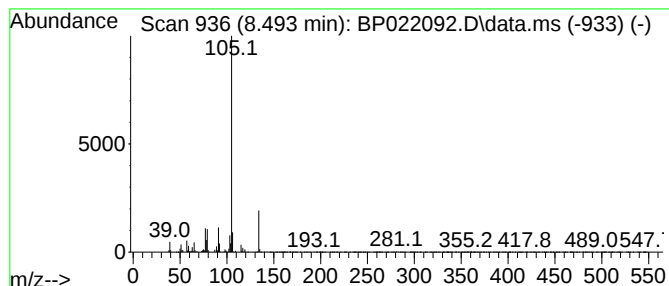
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Benzene, 1-methyl-4-propyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.493	4.57 ng/ul	222389	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	91
2			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
3			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	90
4			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	87
5			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
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Misc :
ALS Vial : 21 Sample Multiplier: 1

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ClientSampleId :
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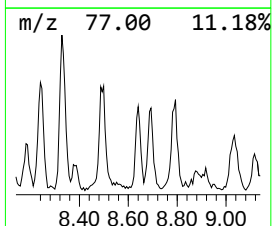
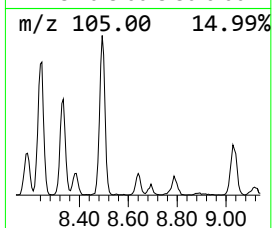
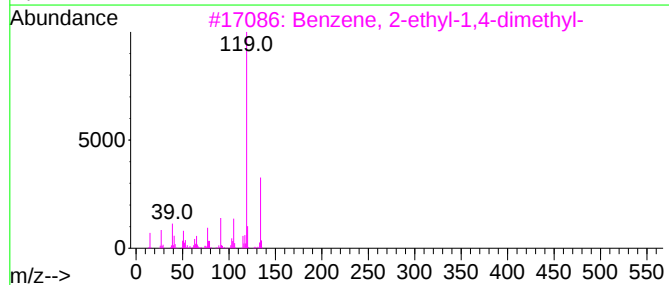
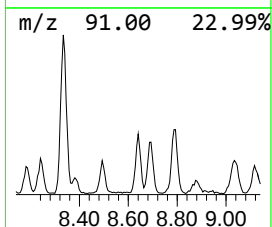
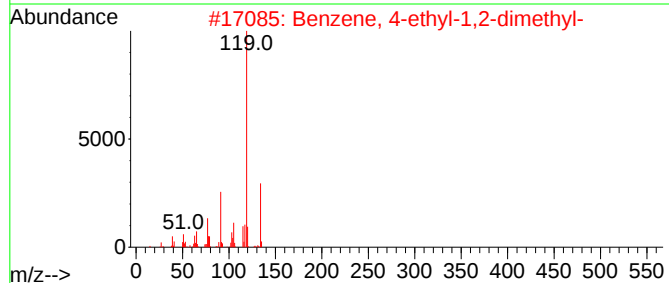
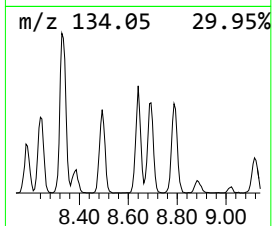
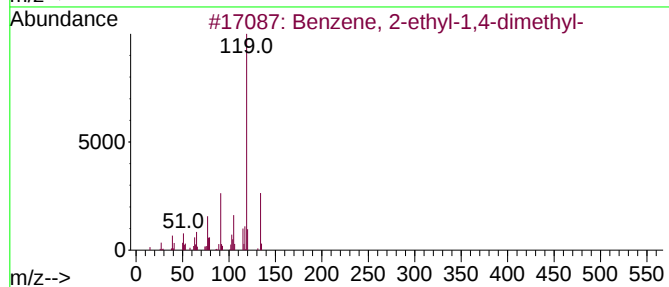
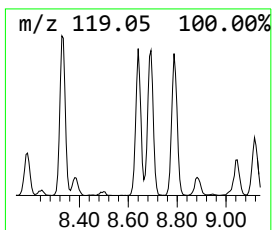
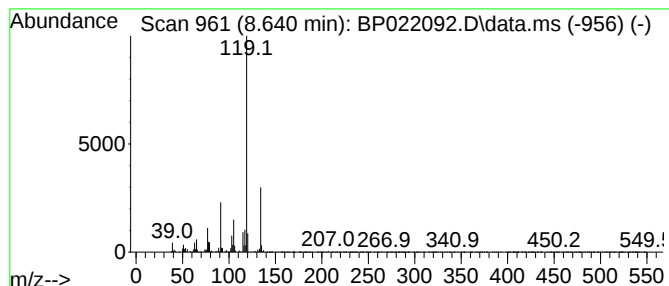
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.640	3.52 ng/ul	171386	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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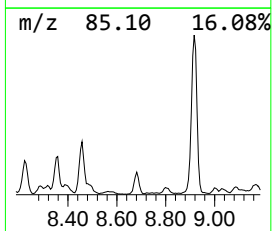
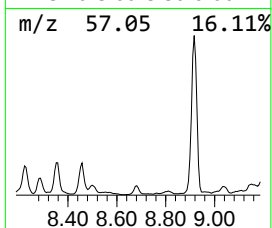
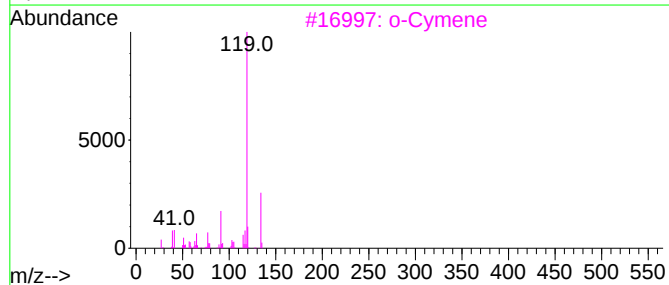
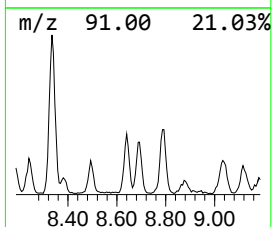
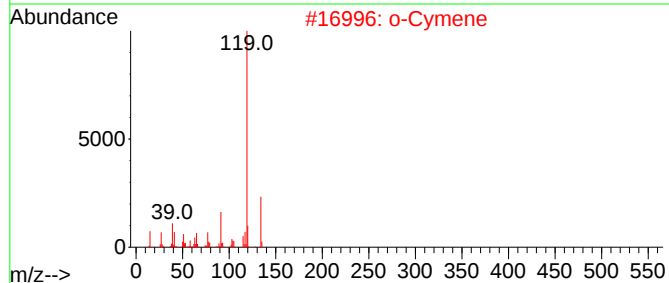
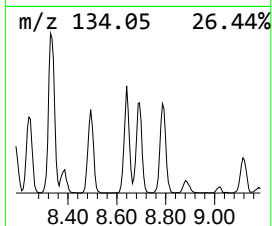
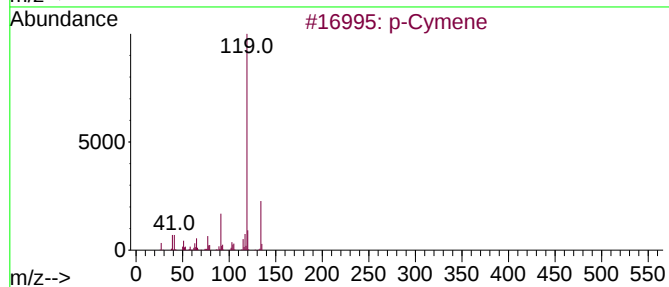
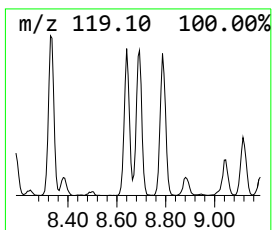
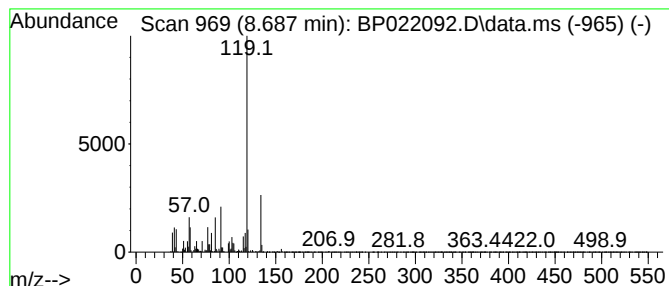
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 p-Cymene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.687	6.08 ng/ul	295972	1,4-Dichlorobenzene-d4	7.699

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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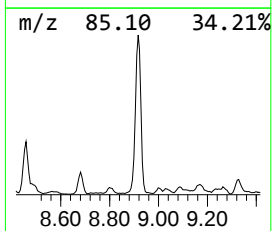
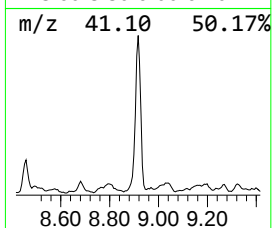
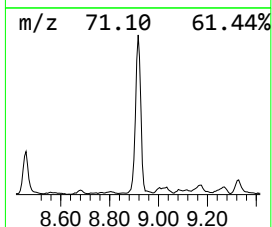
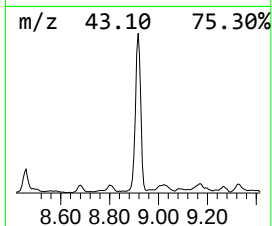
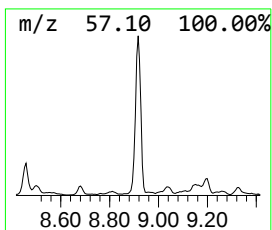
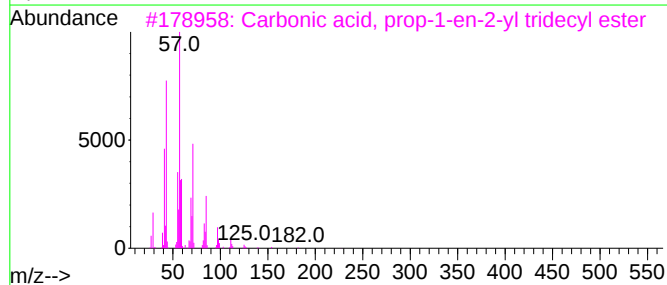
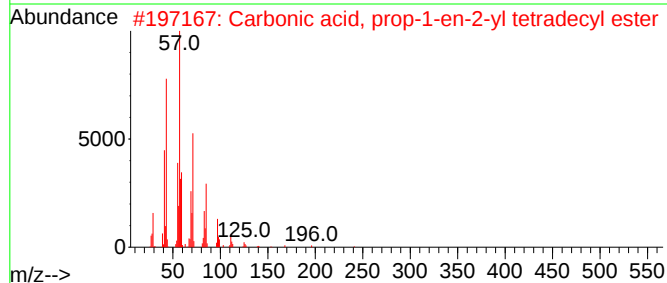
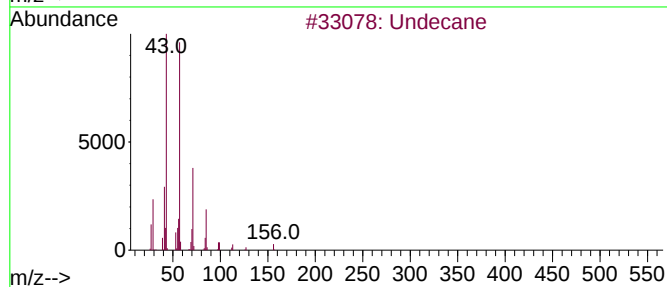
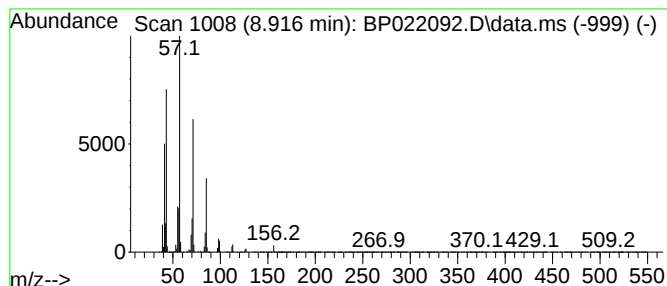
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.916	29.60 ng/ul	1441230	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	91
2			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	86
3			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	83
4			Tridecane	184	C13H28	000629-50-5	80
5			Hexadecane	226	C16H34	000544-76-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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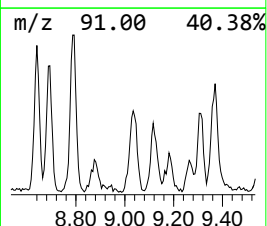
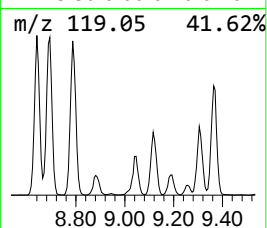
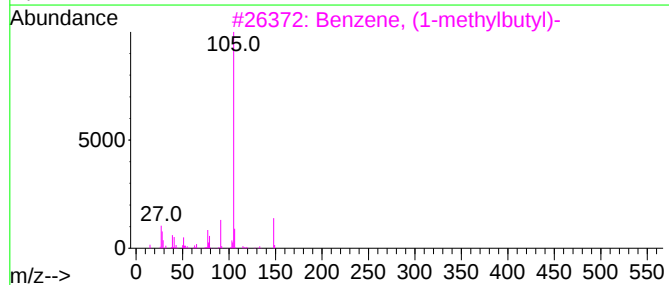
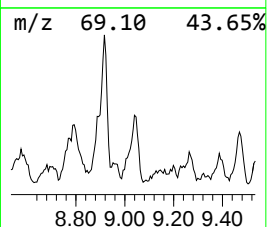
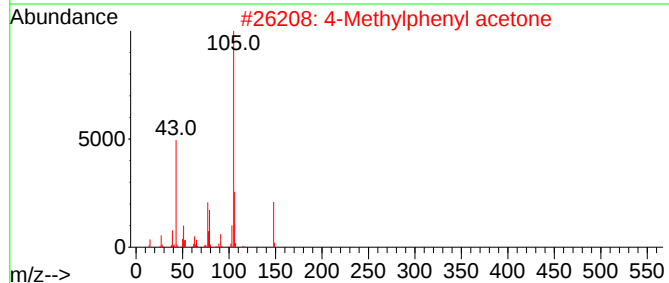
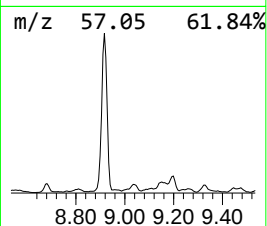
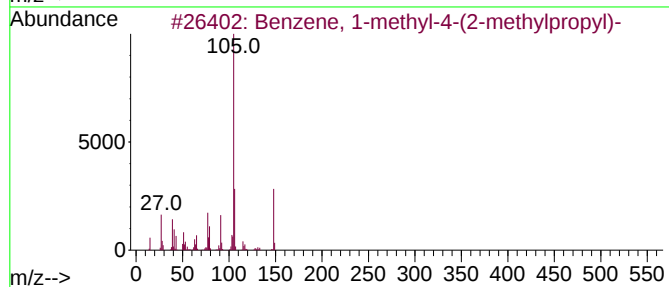
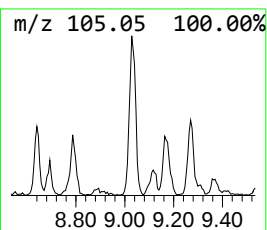
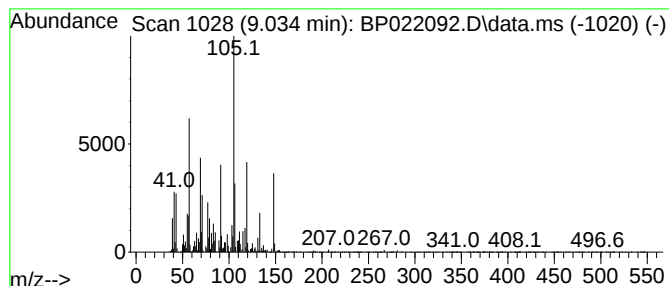
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 unknown-02 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.034	6.89 ng/ul	335421	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	49
2			4-Methylphenyl acetone	148	C10H12O	002096-86-8	45
3			Benzene, (1-methylbutyl)-	148	C11H16	002719-52-0	42
4			3-Buten-2-ol, 4-phenyl-	148	C10H12O	017488-65-2	38
5			2-Methylindan-2-ol	148	C10H12O	033223-84-6	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

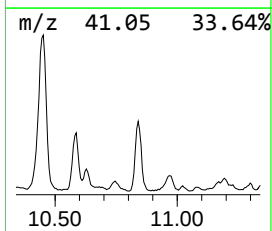
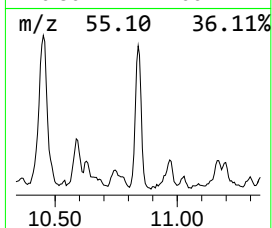
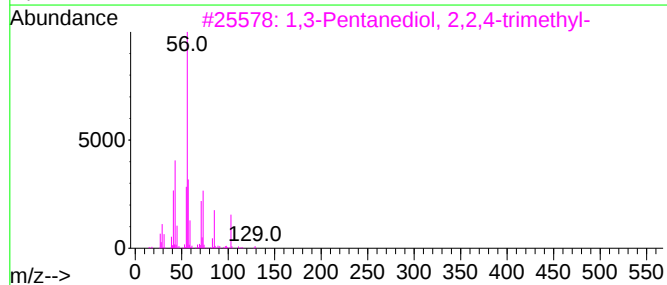
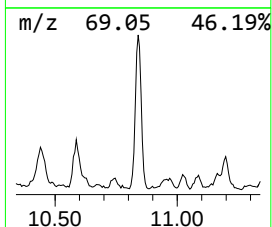
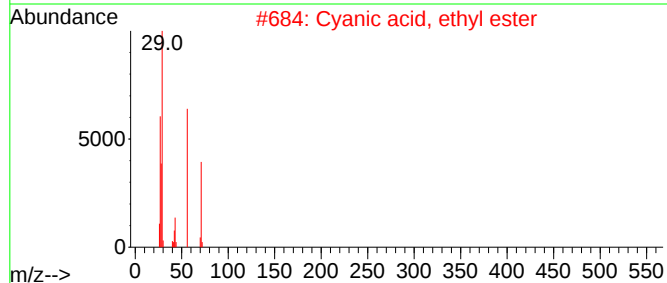
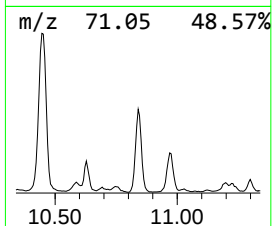
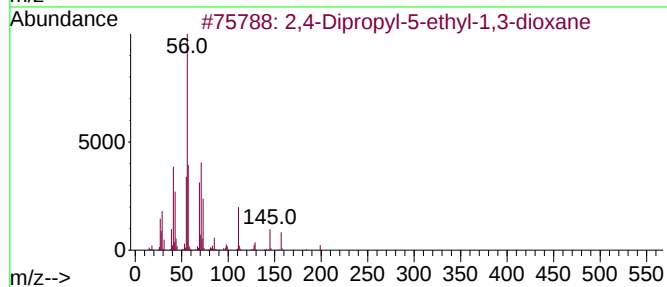
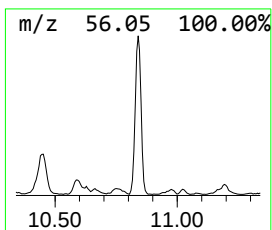
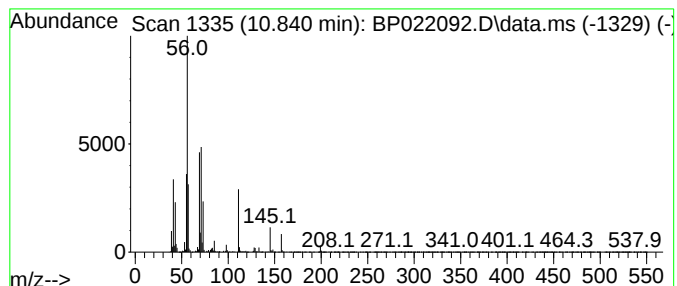
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 2,4-Dipropyl-5-ethyl-1,3-di... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.840	5.72 ng/ul	863178	Naphthalene-d8	10.463	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	64
2	Cyanic acid, ethyl ester	71	C3H5NO	000627-48-5	52
3	1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	28
4	Cyclobutane, 1,2,3,4-tetramethyl-	112	C8H16	069531-57-3	27
5	Tridecane, 7-methylene-	196	C14H28	019780-80-4	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
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Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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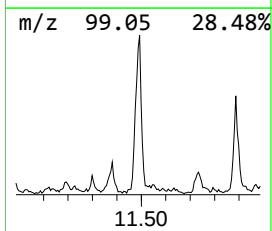
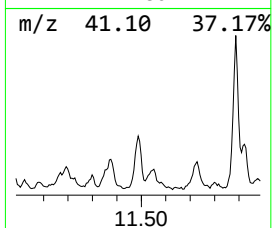
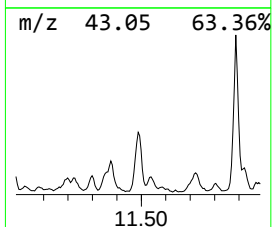
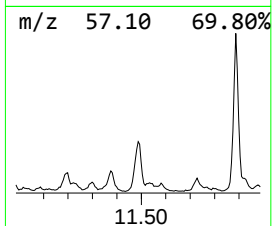
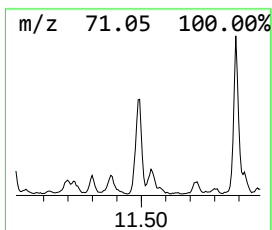
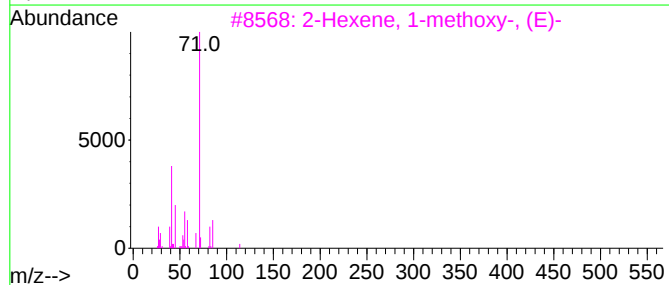
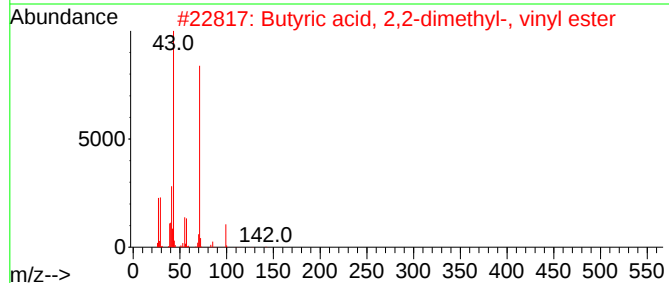
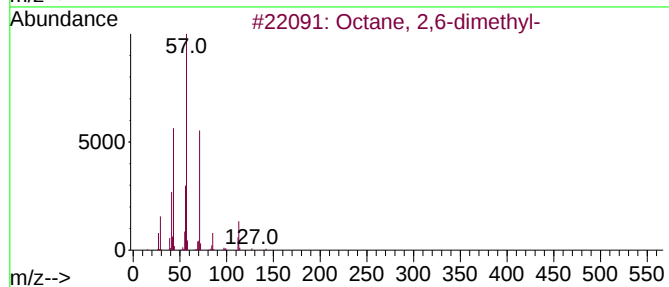
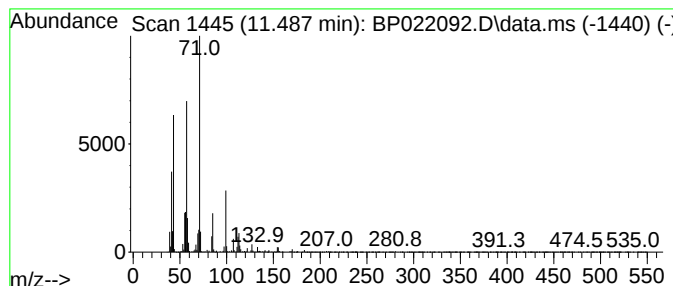
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 56

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.487	2.22 ng/ul	334046	Naphthalene-d8	10.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
2			Butyric acid, 2,2-dimethyl-, vin...	142	C8H14O2	013170-00-8	58
3			2-Hexene, 1-methoxy-, (E)-	114	C7H14O	056052-83-6	53
4			Oxalic acid, neopentyl nonyl ester	286	C16H30O4	1000309-73-3	53
5			Octane, 3-ethyl-	142	C10H22	005881-17-4	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
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Misc :
ALS Vial : 21 Sample Multiplier: 1

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ClientSampleId :
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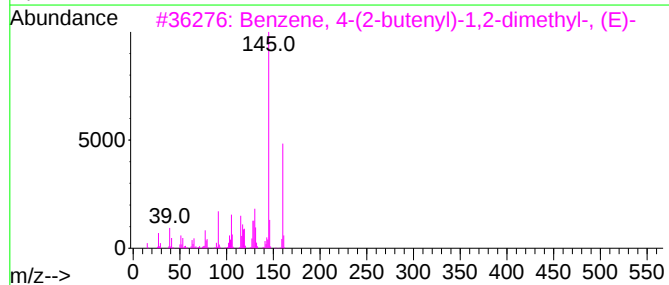
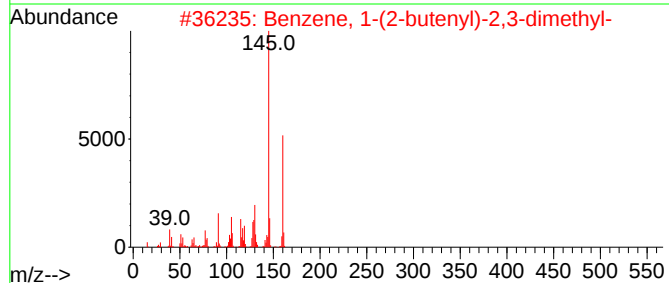
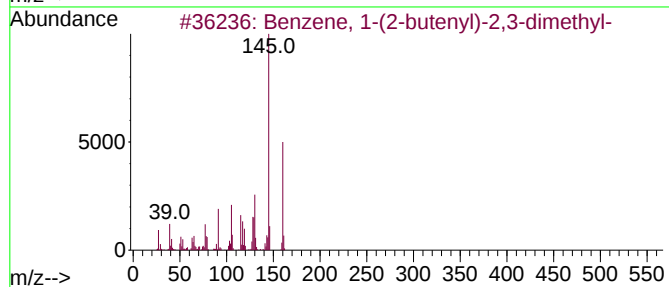
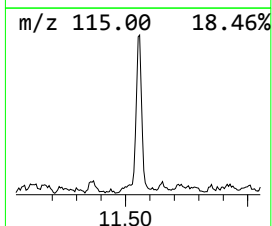
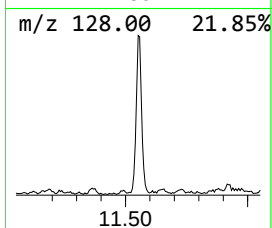
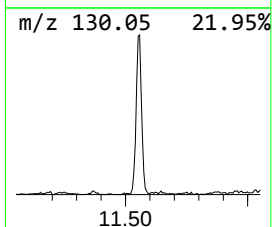
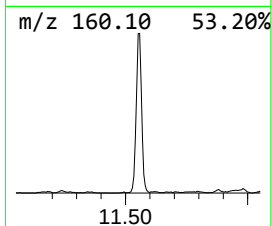
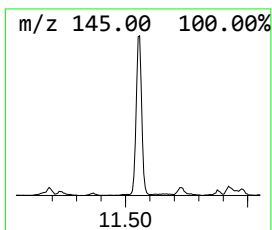
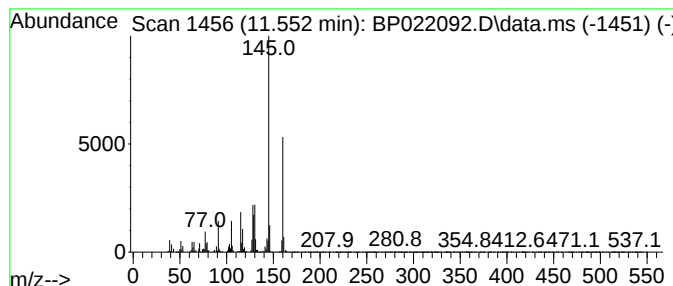
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 Benzene, 1-(2-butenyl)-2,3-... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.552	4.04 ng/ul	608520	Naphthalene-d8	10.463

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	94
2		Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	94
3		Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	94
4		Benzene, 1,2,4-trimethyl-5-(1-me...	160	C12H16	054340-84-0	91
5		Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
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Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

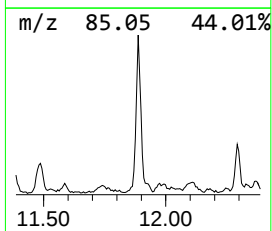
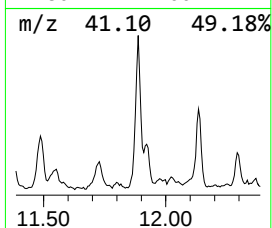
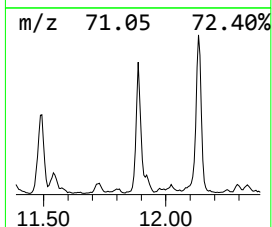
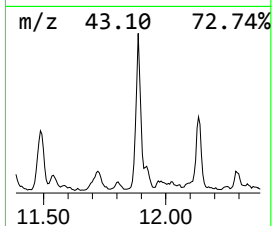
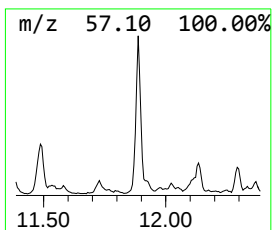
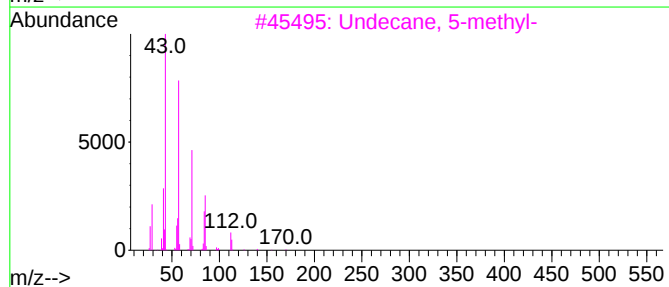
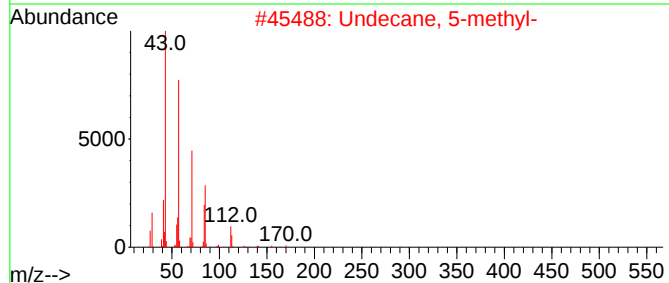
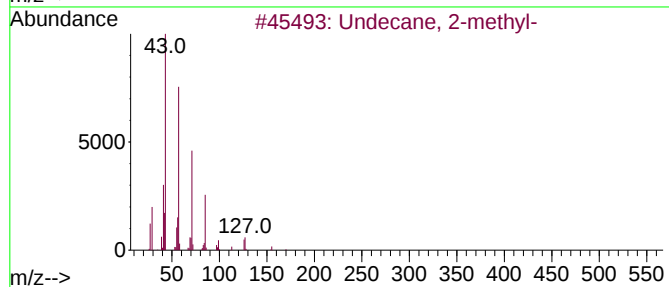
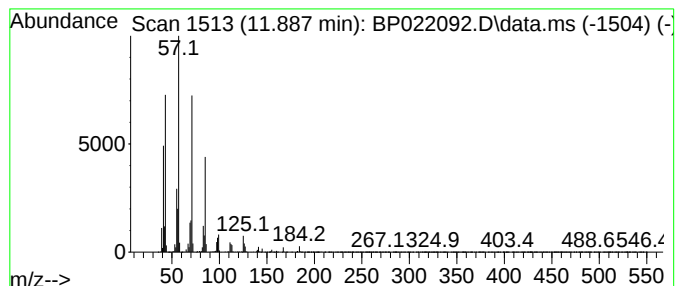
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.887	5.08 ng/ul	766665	Naphthalene-d8	10.463	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 2-methyl-	170	C12H26	007045-71-8	87
2	Undecane, 5-methyl-	170	C12H26	001632-70-8	83
3	Undecane, 5-methyl-	170	C12H26	001632-70-8	80
4	Nonadecane	268	C19H40	000629-92-5	80
5	Tridecane	184	C13H28	000629-50-5	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

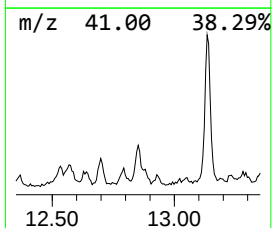
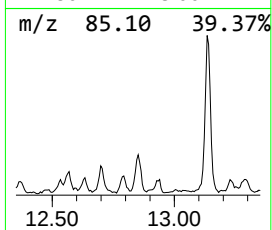
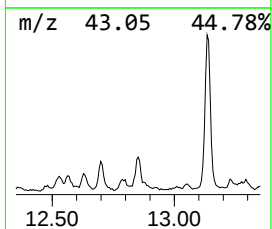
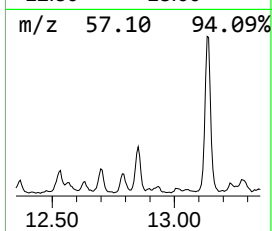
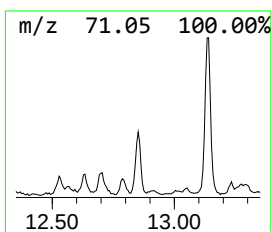
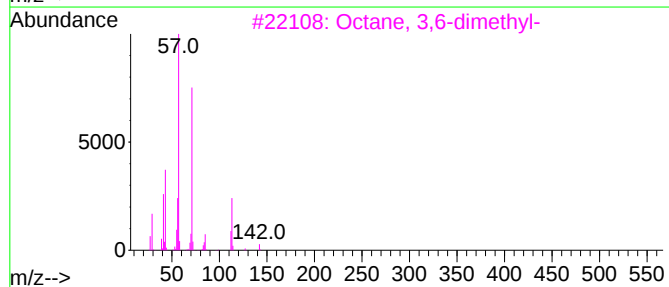
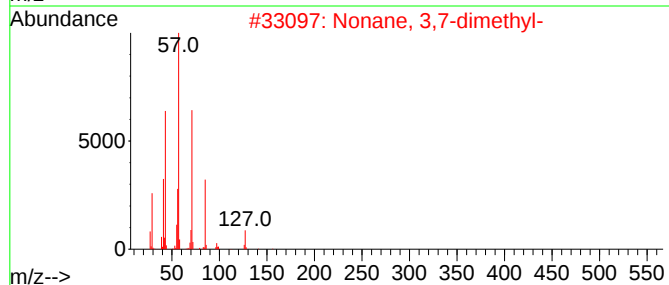
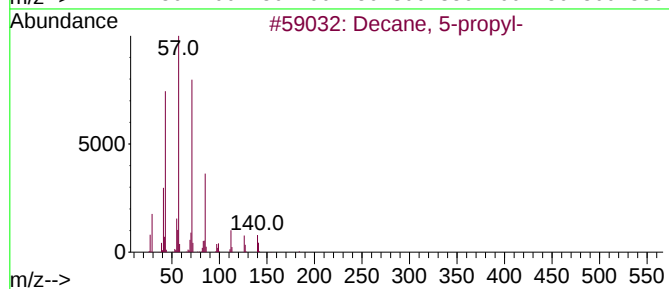
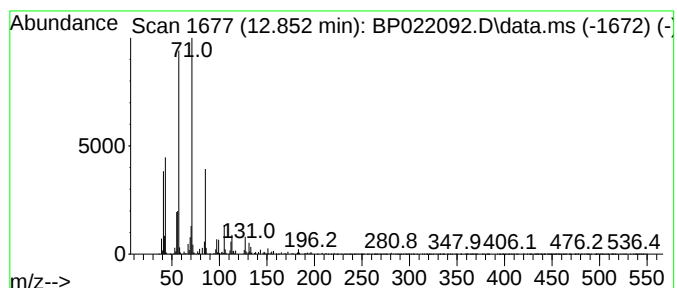
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 57

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.852	2.18 ng/ul	171563	Acenaphthene-d10	14.316	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 5-propyl-	184	C13H28	017312-62-8	74
2	Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	68
3	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	64
4	Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	64
5	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

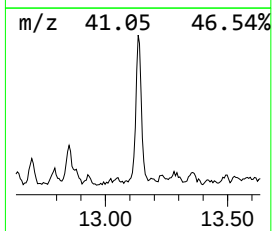
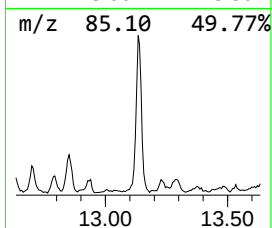
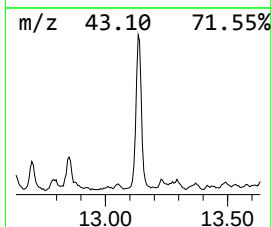
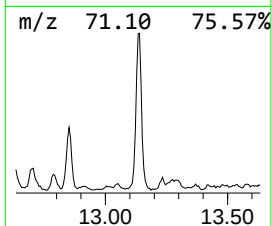
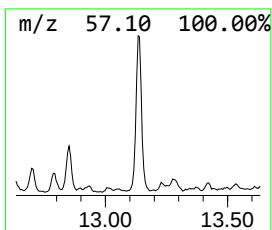
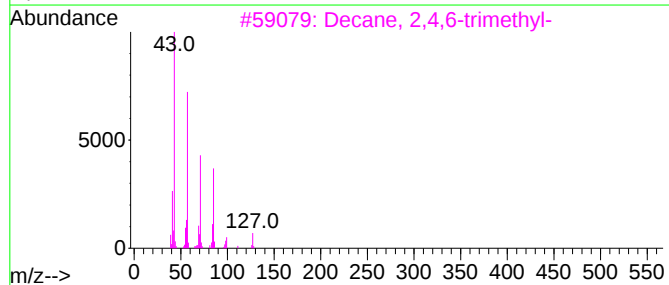
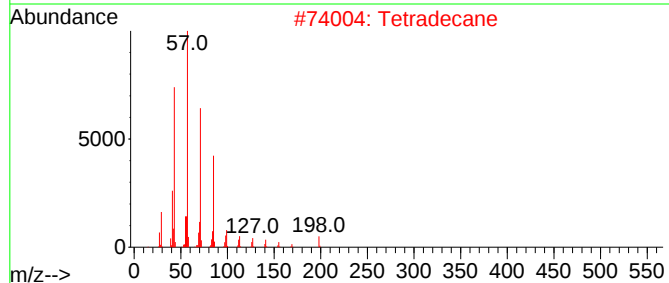
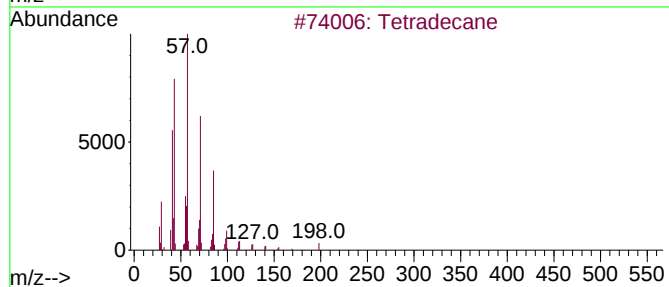
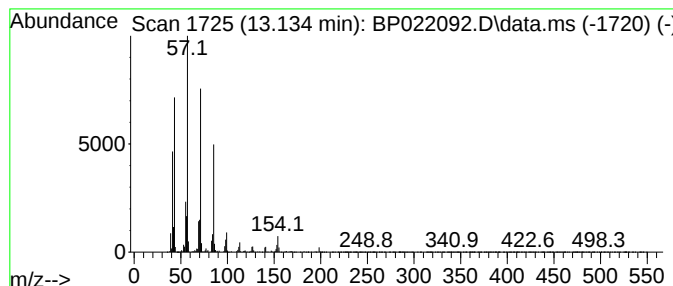
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.134	7.95 ng/ul	625958	Acenaphthene-d10	14.316	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	94
2	Tetradecane	198	C14H30	000629-59-4	90
3	Decane, 2,4,6-trimethyl-	184	C13H28	062108-27-4	86
4	1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	80
5	Hexadecane	226	C16H34	000544-76-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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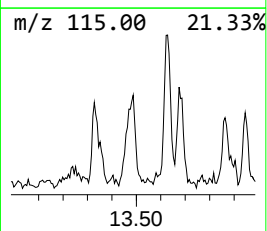
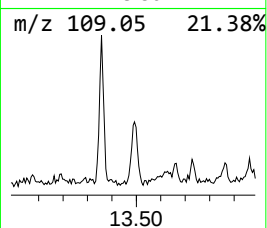
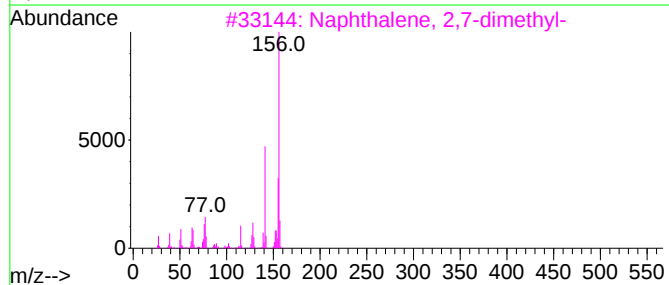
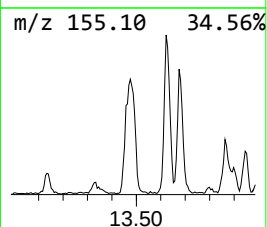
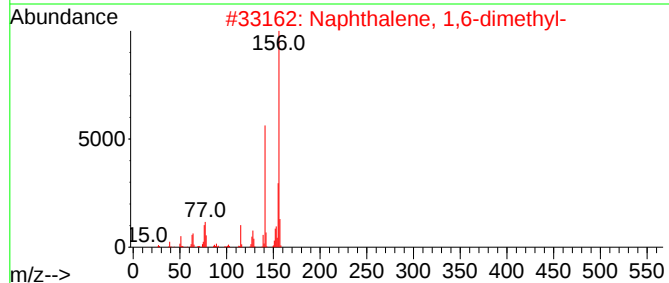
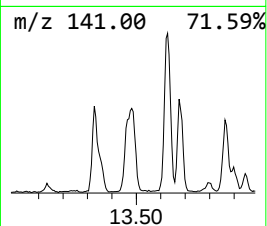
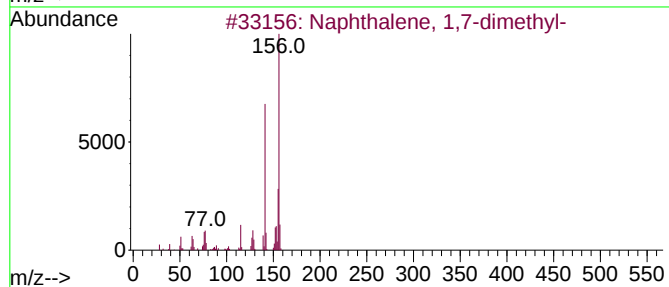
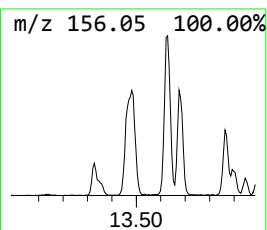
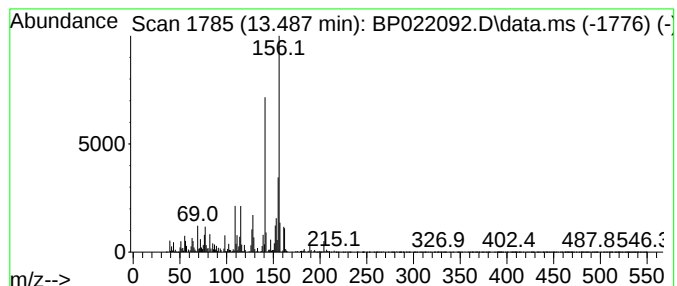
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Quant Title : SVOA CALIBRATION

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

Peak Number 34 Naphthalene, 1,7-dimethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.487	4.43 ng/ul	348986	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
5			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
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Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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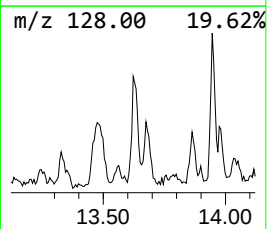
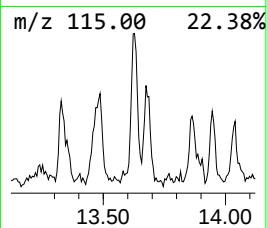
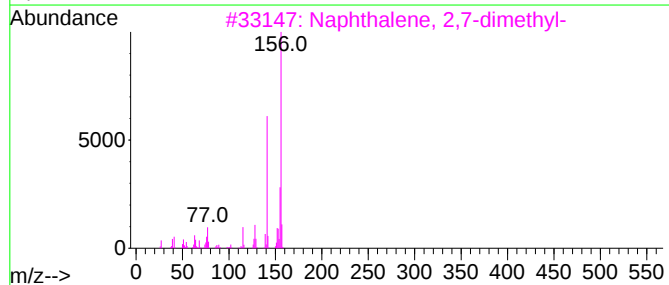
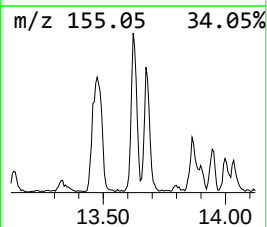
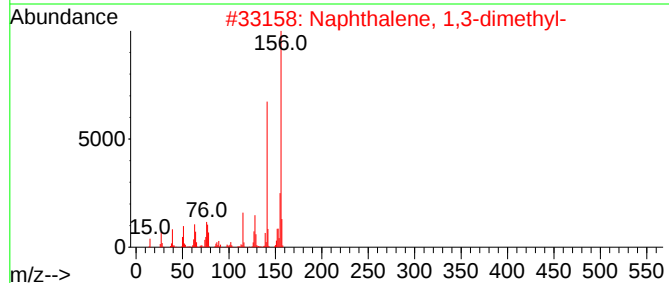
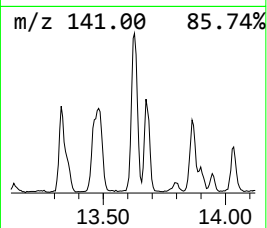
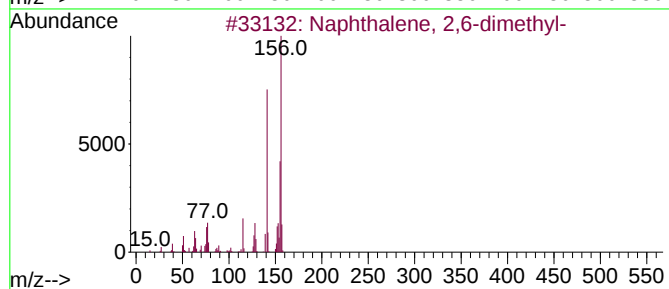
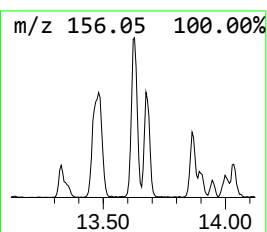
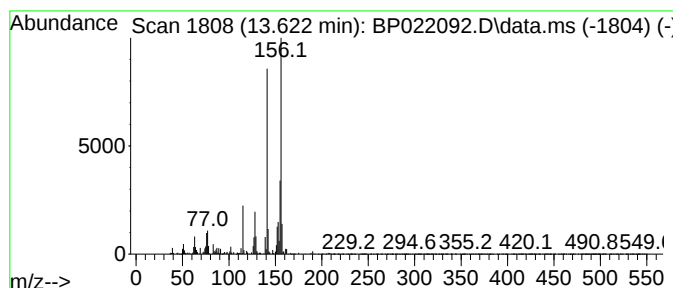
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Quant Title : SVOA CALIBRATION

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

Peak Number 35 Naphthalene, 2,6-dimethyl- Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.622	3.66 ng/ul	287869	Acenaphthene-d10	14.316

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
2		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
3		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
4		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
5		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC63

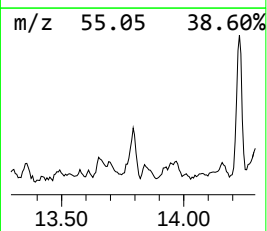
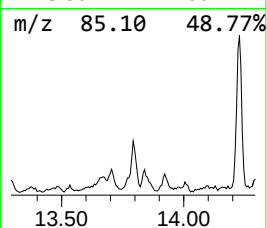
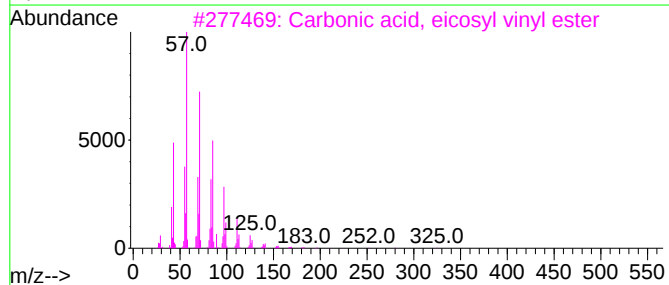
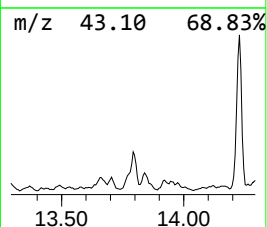
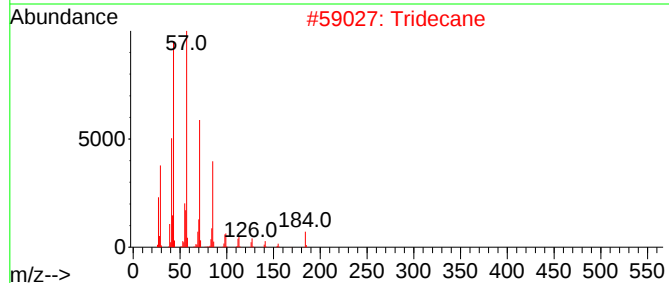
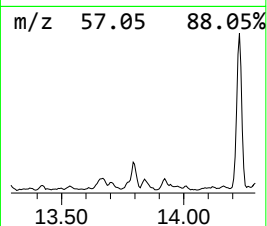
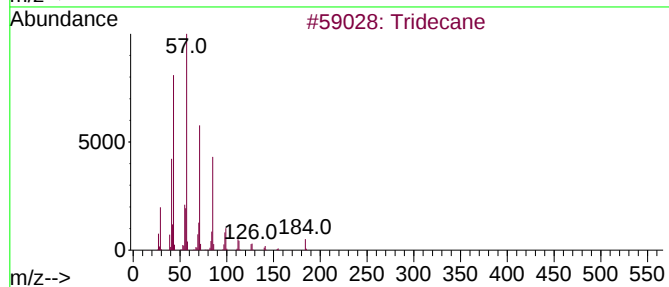
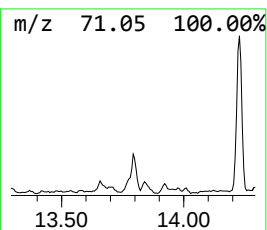
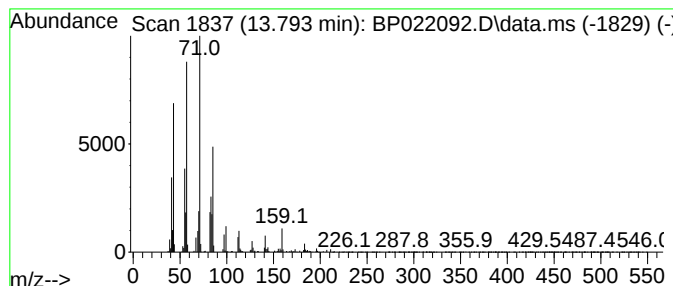
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Quant Title : SVOA CALIBRATION

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

Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.793	3.45 ng/ul	271919	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	64
2			Tridecane	184	C13H28	000629-50-5	64
3			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	58
4			Tridecane, 7-propyl-	226	C16H34	055045-09-5	55
5			Tridecane, 6-propyl-	226	C16H34	055045-10-8	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

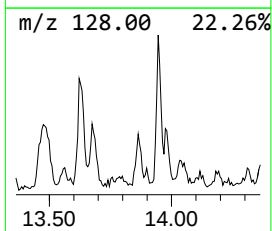
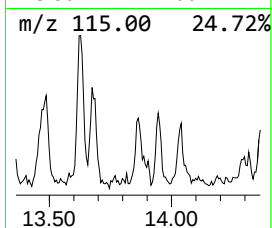
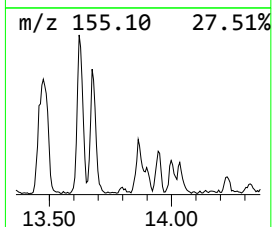
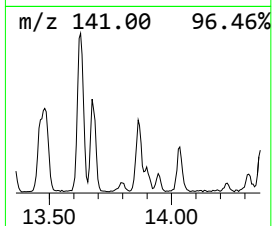
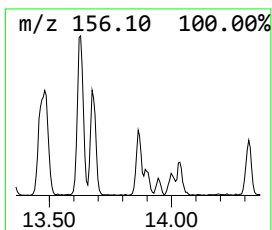
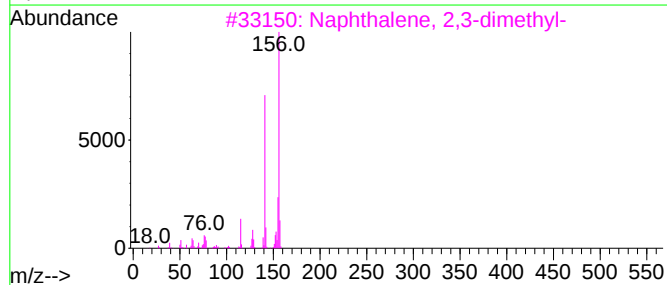
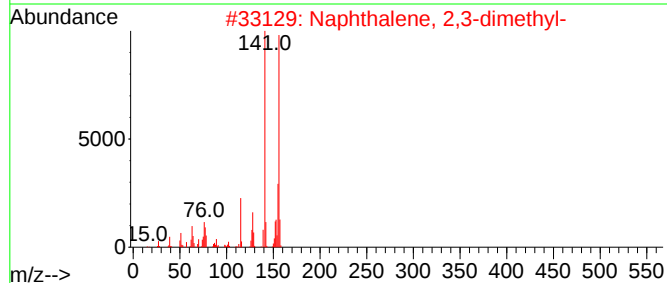
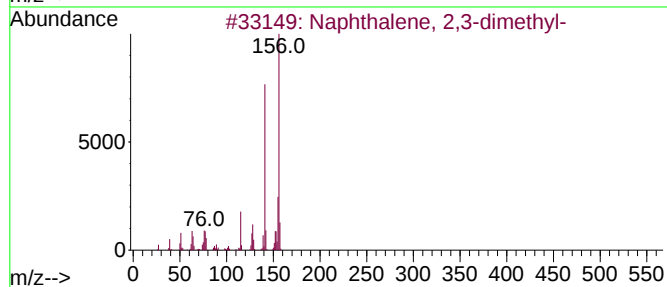
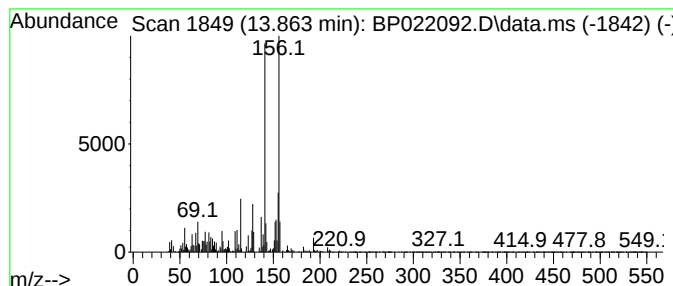
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 37 Naphthalene, 2,3-dimethyl- Concentration Rank 53

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.863	2.68 ng/ul	211442	Acenaphthene-d10	14.316	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	94
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	93
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	93
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	93
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC63

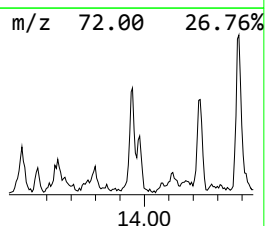
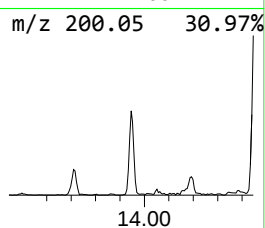
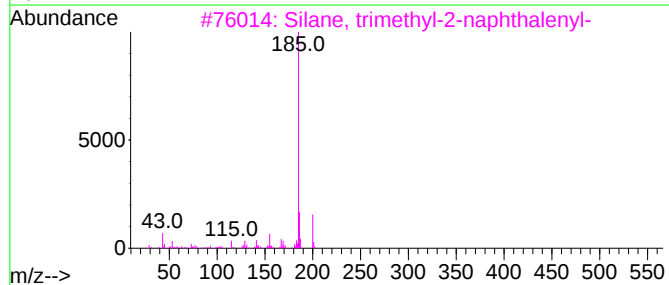
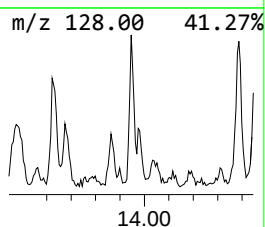
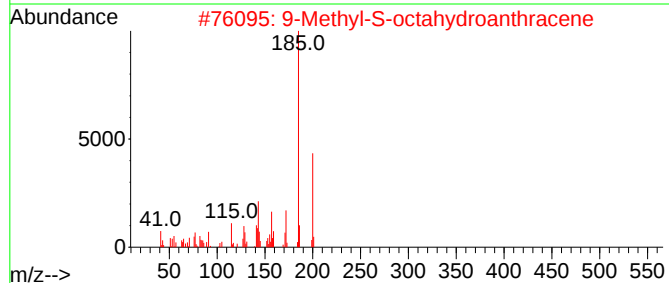
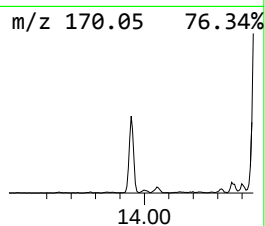
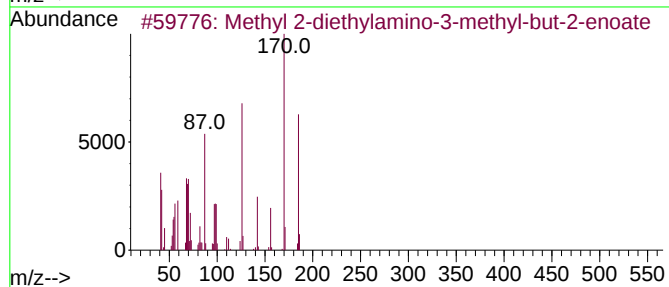
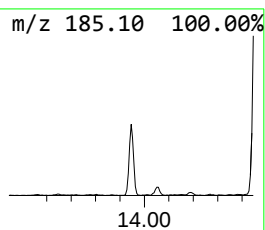
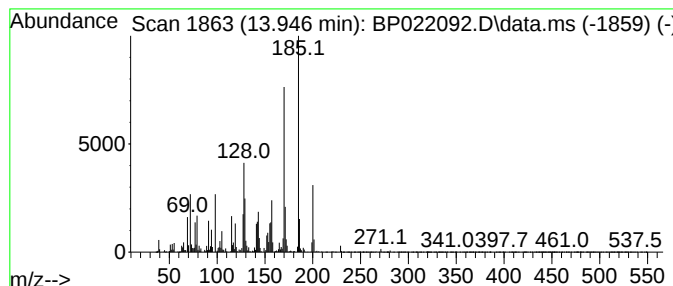
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 unknown-03 Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.946	2.90 ng/ul	228621	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 2-diethylamino-3-methyl-b...	185	C10H19NO2	075122-81-5	49
2			9-Methyl-S-octahydroanthracene	200	C15H20	004948-51-0	42
3			Silane, trimethyl-2-naphthalenyl-	200	C13H16Si	018052-85-2	42
4			1,2,3,4,5,6,7,8-Octahydro-1-meth...	200	C15H20	1000080-19-4	42
5			.gamma.-Dehydro-ar-himachalene	200	C15H20	051766-65-5	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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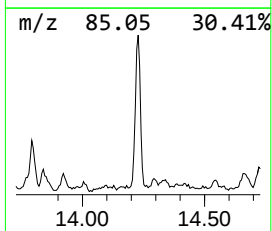
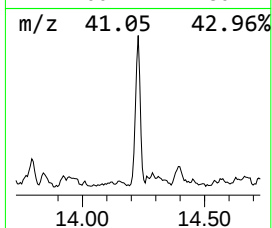
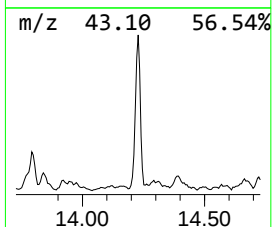
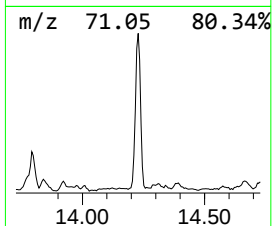
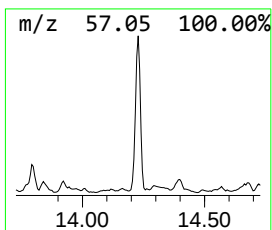
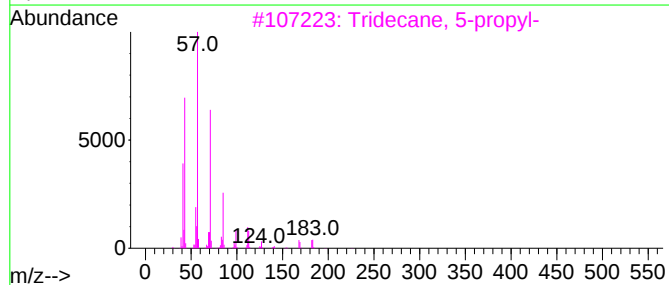
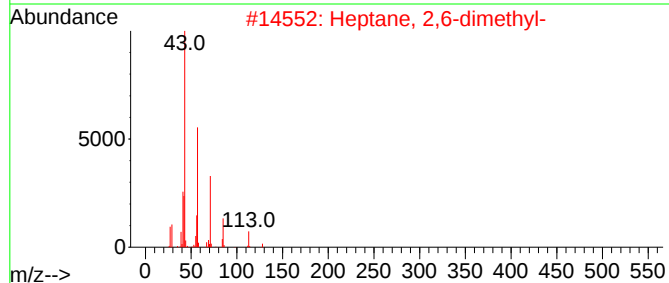
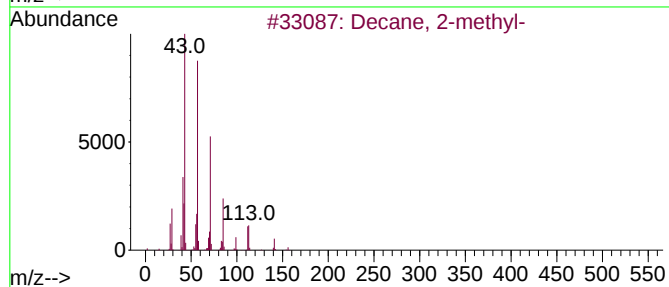
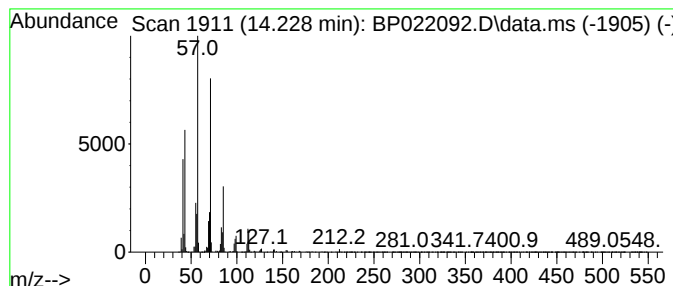
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.228	11.28 ng/ul	888690	Acenaphthene-d10	14.316

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 2-methyl-	156	C11H24	006975-98-0	83
2		Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	80
3		Tridecane, 5-propyl-	226	C16H34	055045-11-9	78
4		Decane, 2-methyl-	156	C11H24	006975-98-0	78
5		Pentadecane, 7-methyl-	226	C16H34	006165-40-8	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC63

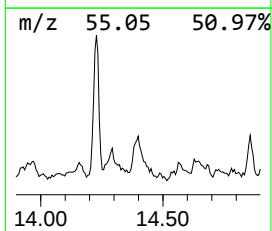
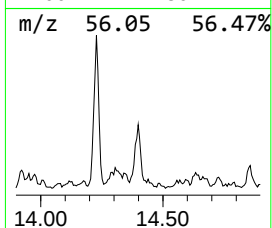
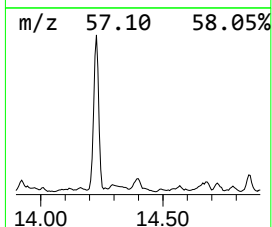
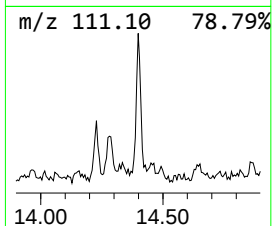
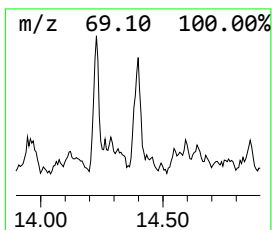
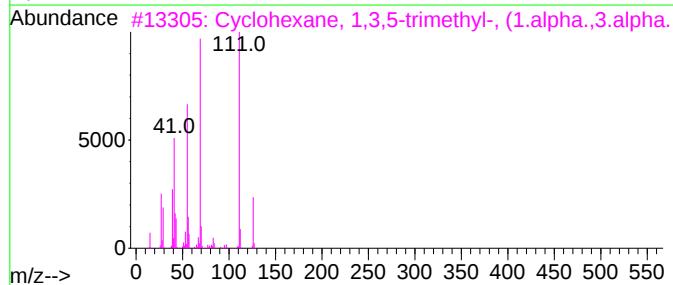
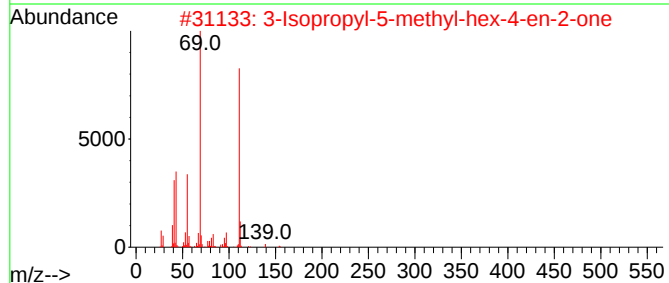
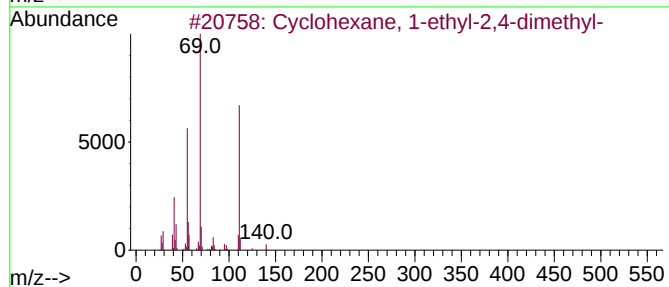
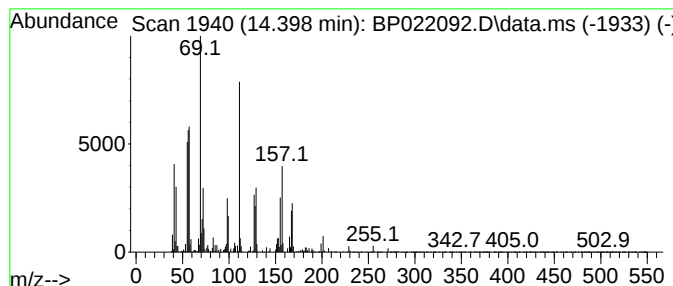
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Quant Title : SVOA CALIBRATION

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

Peak Number 40 (DEL) Alkane: Cyclic14.399 Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.399	3.13 ng/ul	246518	Acenaphthene-d10	14.316

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 1-ethyl-2,4-dimethyl-	140	C10H20	061142-69-6	41
2		3-Isopropyl-5-methyl-hex-4-en-2-one	154	C10H18O	077142-85-9	38
3		Cyclohexane, 1,3,5-trimethyl-, (...)	126	C9H18	001795-27-3	35
4		Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	35
5		Cyclohexane, 1-ethyl-1,3-dimethy...	140	C10H20	062238-31-7	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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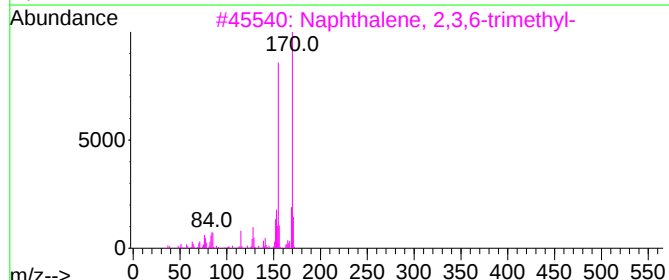
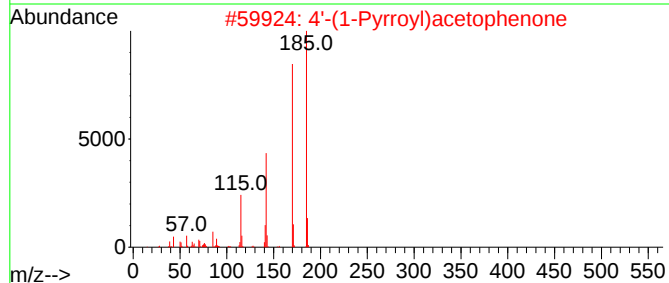
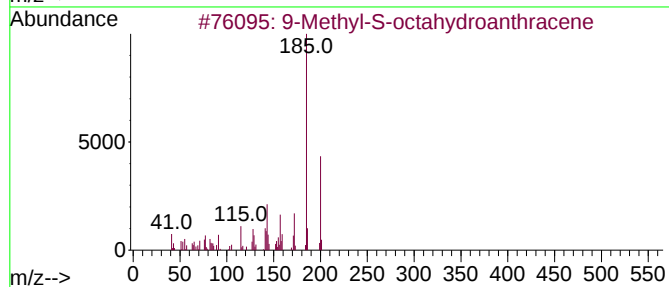
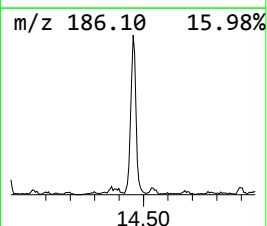
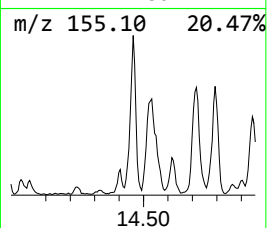
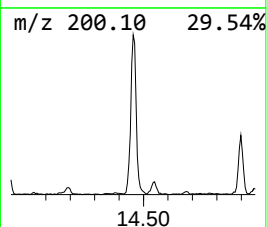
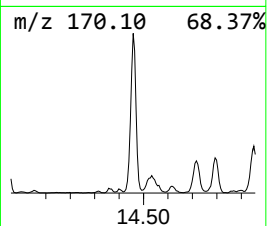
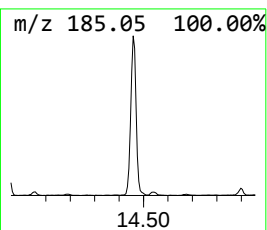
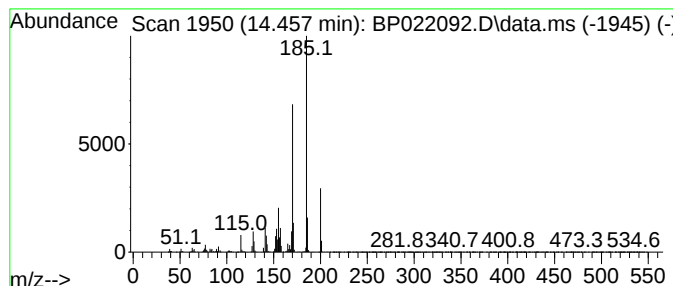
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Quant Title : SVOA CALIBRATION

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

Peak Number 41 9-Methyl-S-octahydroanthracene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.457	9.68 ng/ul	762451	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Methyl-S-octahydroanthracene	200	C15H20	004948-51-0	60
2			4'-(1-Pyrrolyl)acetophenone	185	C12H11NO	022106-37-2	52
3			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	51
4			1H-Pyrazole, 5-(t-butyl)-3-phenyl-	200	C13H16N2	1000261-34-8	46
5			Ethanone, 1-(6-methoxy-1-naphtha...	200	C13H12O2	058149-89-6	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

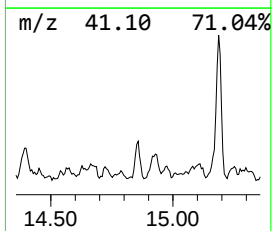
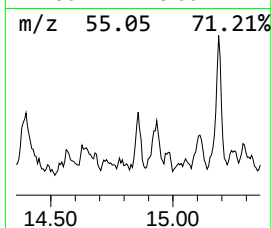
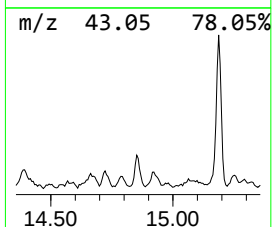
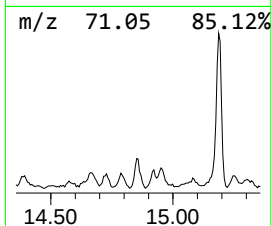
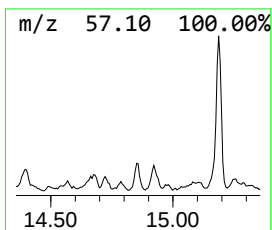
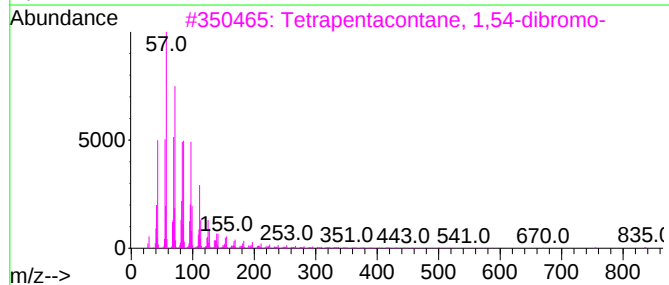
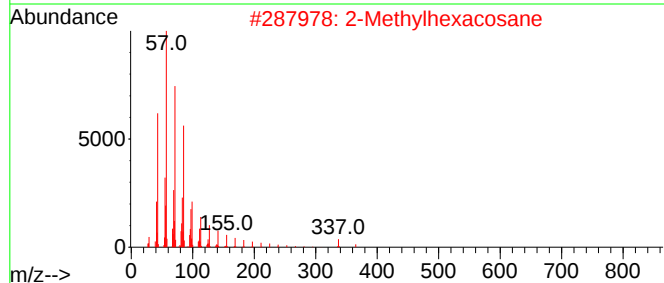
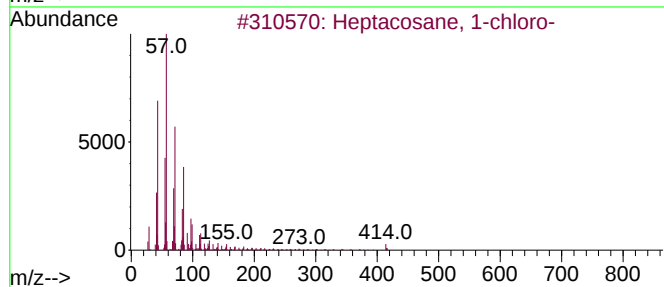
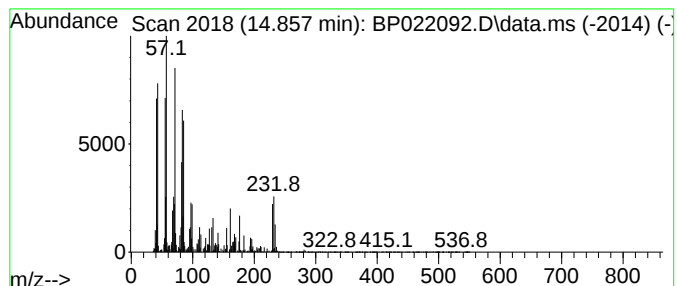
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 Heptacosane, 1-chloro- Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.857	2.79 ng/ul	219635	Acenaphthene-d10	14.316	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	55
2	2-Methylhexacosane	380	C27H56	001561-02-0	53
3	Tetrapentacontane, 1,54-dibromo-	915	C54H108Br2	1000156-09-4	52
4	Hexatriacontane	507	C36H74	000630-06-8	49
5	Triacontane, 1-bromo-	500	C30H61Br	004209-22-7	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC63

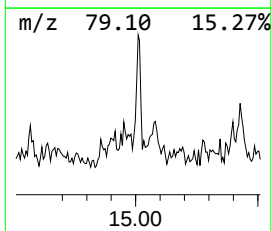
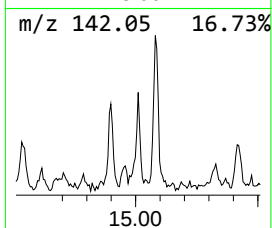
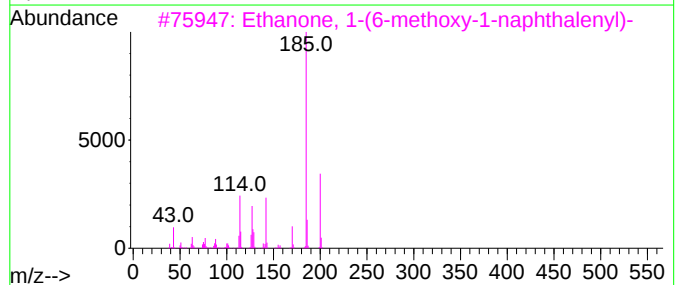
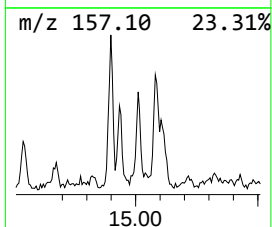
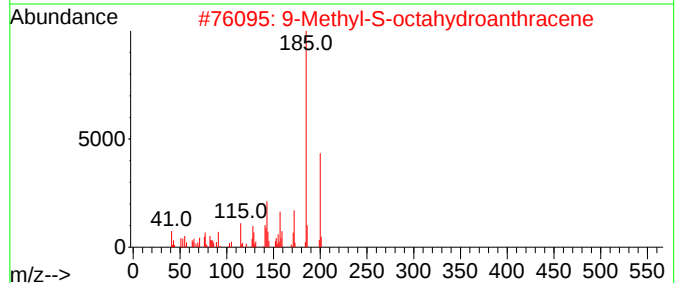
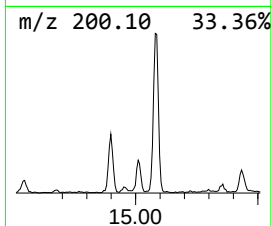
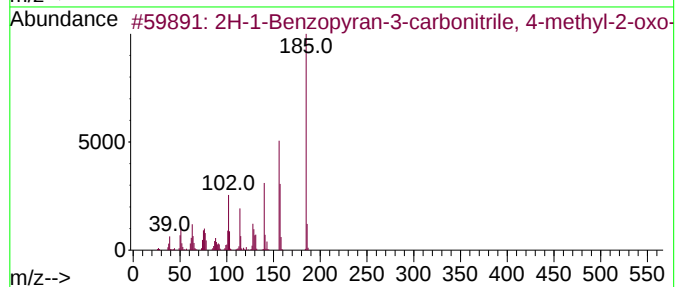
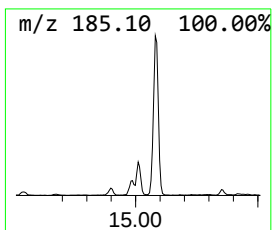
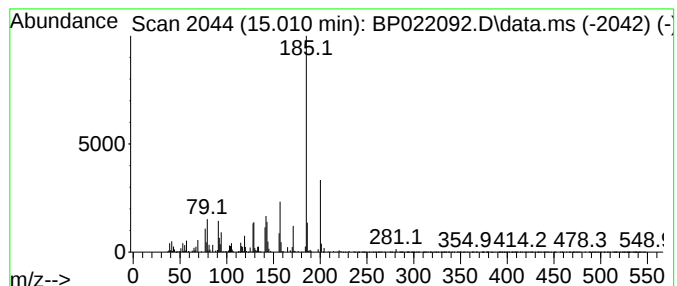
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Quant Title : SVOA CALIBRATION

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

Peak Number 44 2H-1-Benzopyran-3-carbonitr... Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.010	2.35 ng/ul	184741	Acenaphthene-d10	14.316

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2H-1-Benzopyran-3-carbonitrile, ...	185	C11H7NO2	024526-69-0	58
2		9-Methyl-S-octahydroanthracene	200	C15H20	004948-51-0	53
3		Ethanone, 1-(6-methoxy-1-naphtha...	200	C13H12O2	058149-89-6	53
4		Cyclopropane, 1-methoxy-2,2-dime...	200	C14H16O	1000271-83-0	53
5		Benzenesulfonamide, 4-ethylamino-	200	C8H12N2O2S	1000452-18-1	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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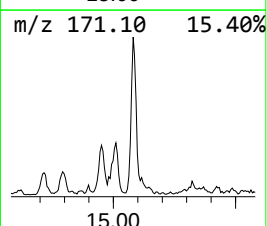
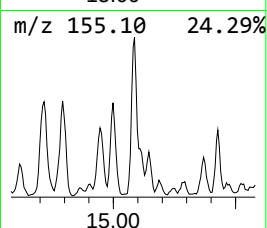
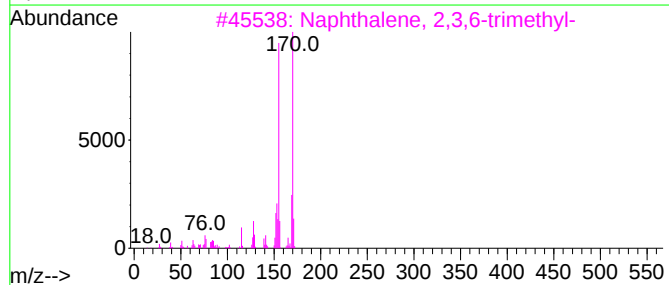
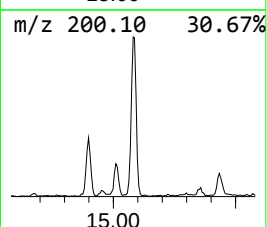
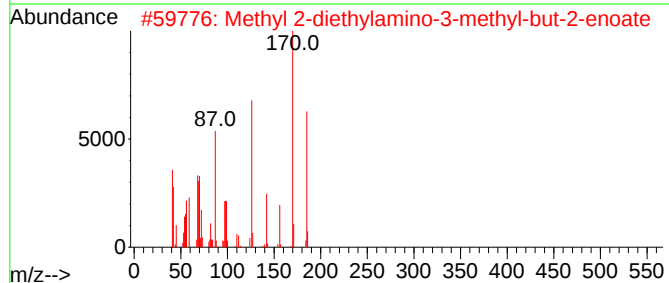
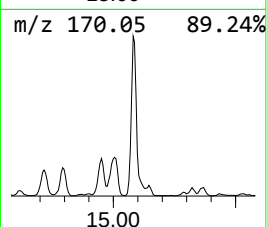
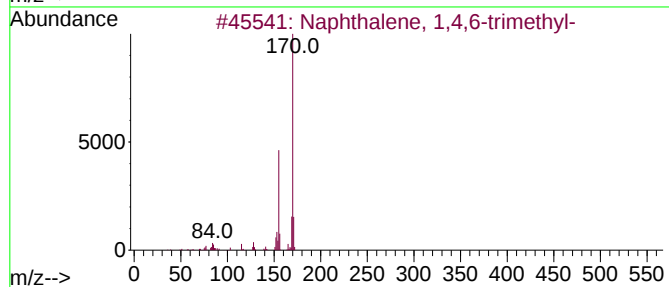
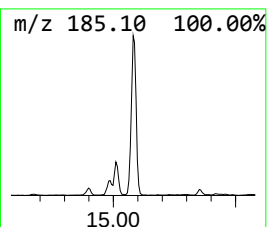
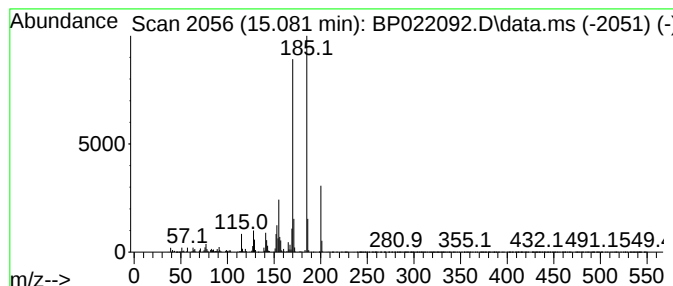
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Quant Title : SVOA CALIBRATION

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

Peak Number 45 Naphthalene, 1,4,6-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.081	11.04 ng/ul	869585	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	55
2			Methyl 2-diethylamino-3-methyl-b...	185	C10H19NO2	075122-81-5	50
3			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	45
4			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	45
5			Ethanone, 1-(6-methoxy-1-naphtha...	200	C13H12O2	058149-89-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

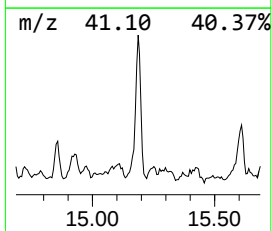
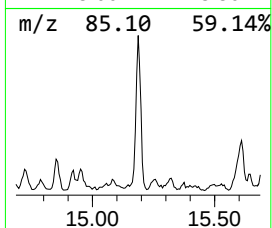
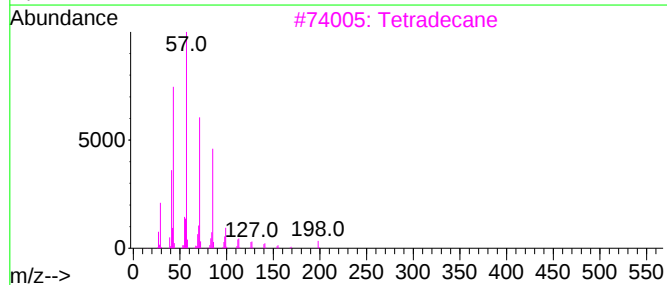
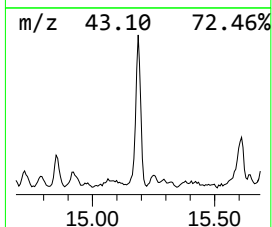
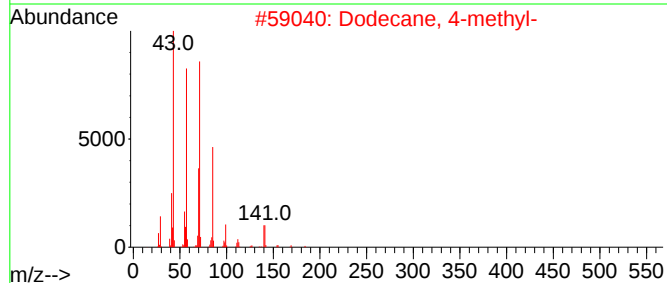
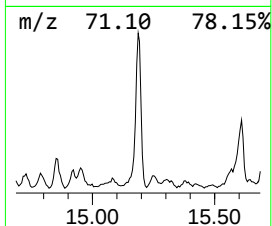
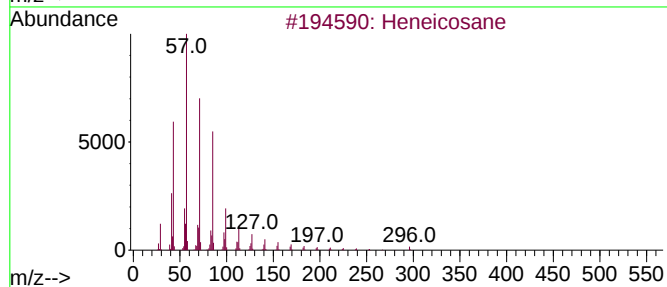
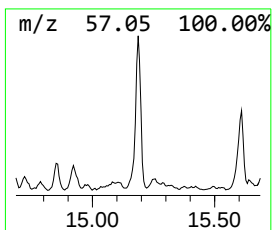
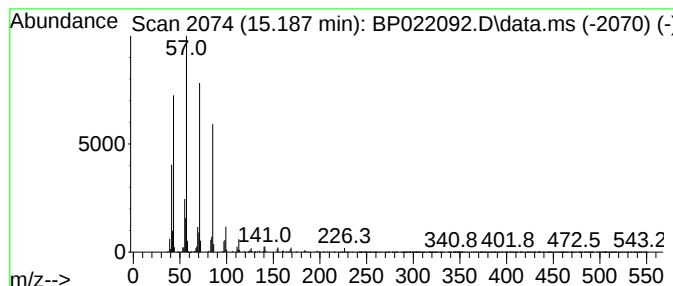
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 46 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.187	7.03 ng/ul	553580	Acenaphthene-d10		14.316	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Dodecane, 4-methyl-		184	C13H28	006117-97-1	90
3	Tetradecane		198	C14H30	000629-59-4	86
4	Eicosane		282	C20H42	000112-95-8	72
5	Pentadecane		212	C15H32	000629-62-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

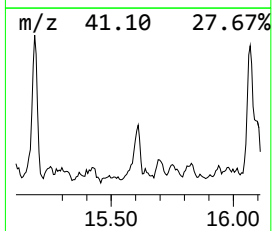
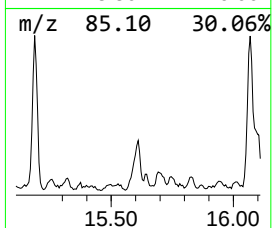
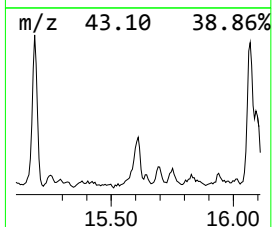
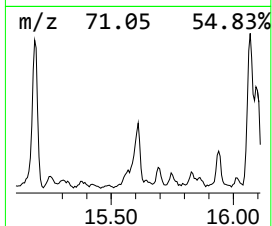
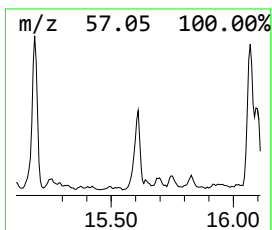
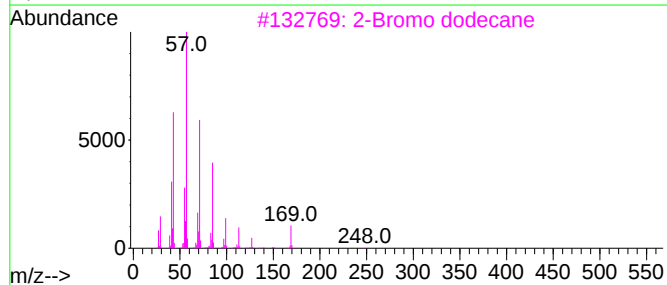
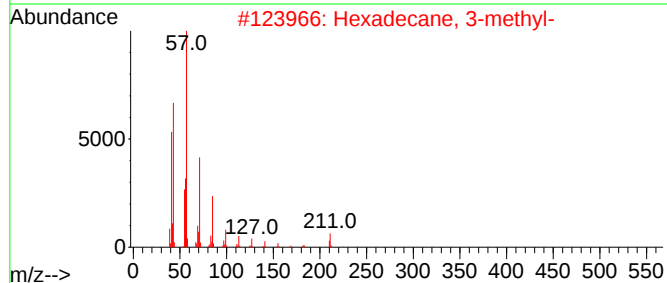
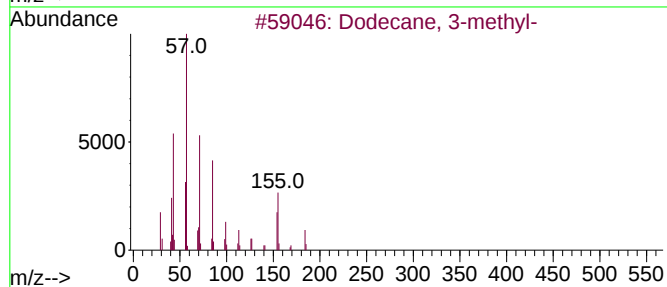
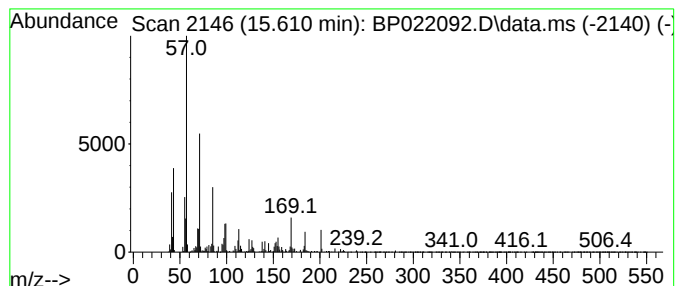
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.610	3.76 ng/ul	296125	Acenaphthene-d10	14.316	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 3-methyl-	184	C13H28	017312-57-1	81
2	Hexadecane, 3-methyl-	240	C17H36	006418-43-5	70
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	64
4	Eicosane, 2-methyl-	296	C21H44	001560-84-5	58
5	Docosane, 11-butyl-	366	C26H54	013475-76-8	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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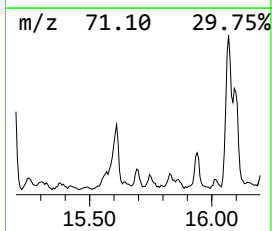
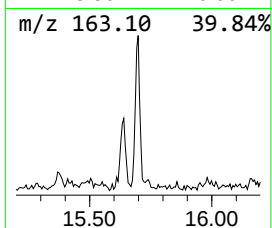
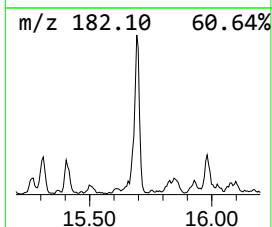
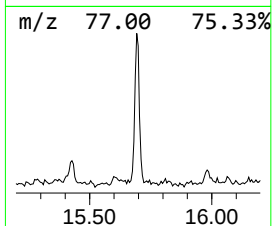
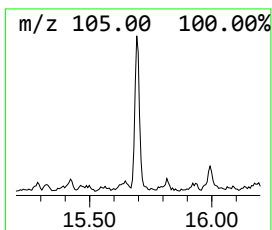
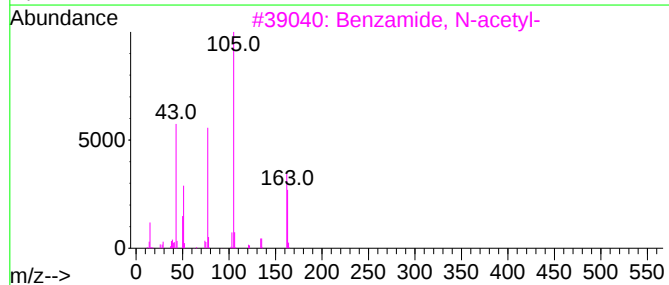
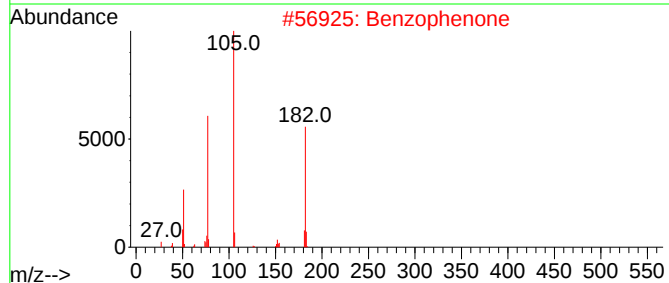
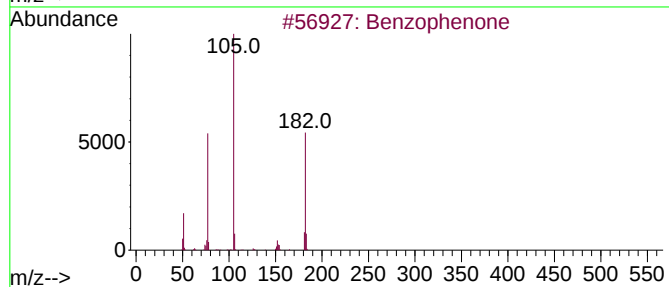
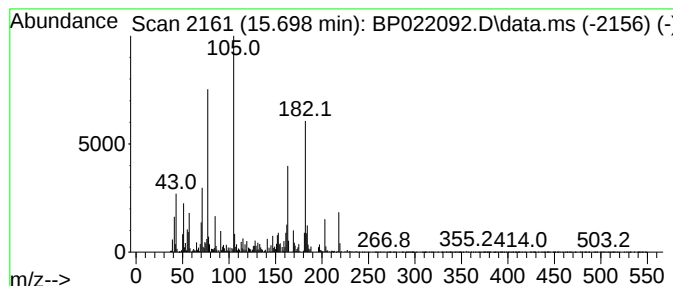
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Quant Title : SVOA CALIBRATION

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

Peak Number 48 Benzophenone Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.698	3.87 ng/ul	304814	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	91
2			Benzophenone	182	C13H10O	000119-61-9	60
3			Benzamide, N-acetyl-	163	C9H9NO2	001575-95-7	49
4			4-Fluoro-5-phenyl-1,2-oxazole	163	C9H6FNO	1000411-55-8	47
5			Azobenzene	182	C12H10N2	000103-33-3	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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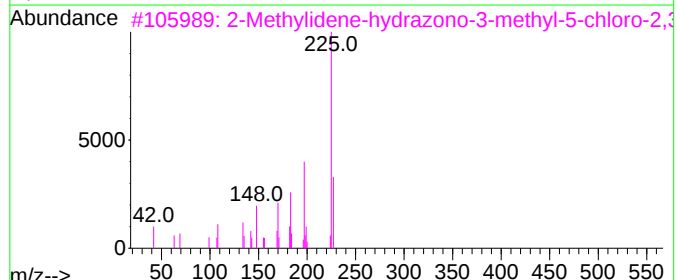
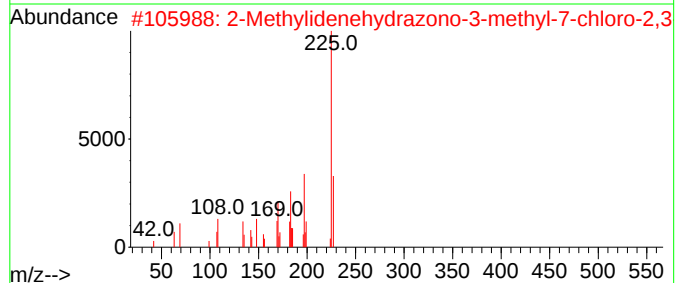
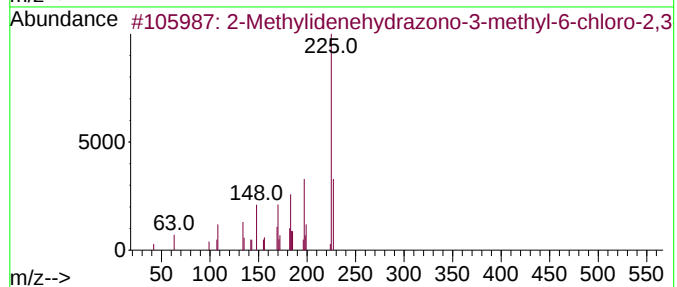
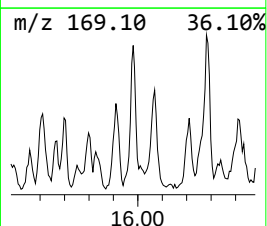
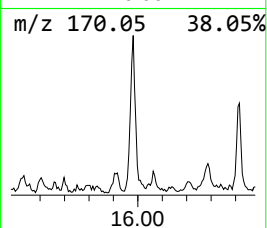
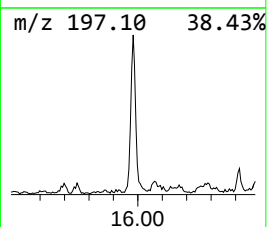
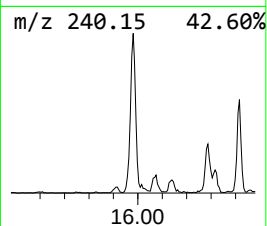
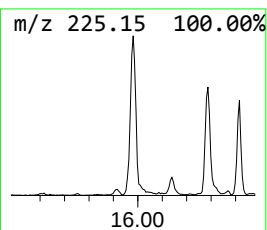
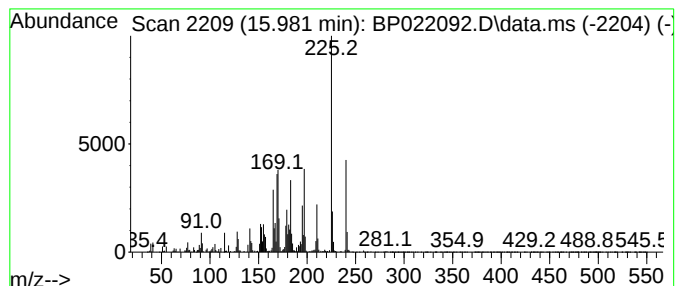
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Quant Title : SVOA CALIBRATION

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

Peak Number 49 unknown-04 Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.981	3.78 ng/ul	359979	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylidenehydrazono-3-methyl-...	225	C9H8ClN3S	1000147-91-5	45		
2	2-Methylidenehydrazono-3-methyl-...	225	C9H8ClN3S	1000147-91-6	43		
3	2-Methylidene-hydrazono-3-methyl...	225	C9H8ClN3S	1000147-91-3	41		
4	(3-Isopropyl-5-methylamino-2,4,6...	225	C14H15N3	014204-00-3	38		
5	6-Aminobenzo(g)quinoxaline-5,10-...	225	C12H7N3O2	072225-24-2	35		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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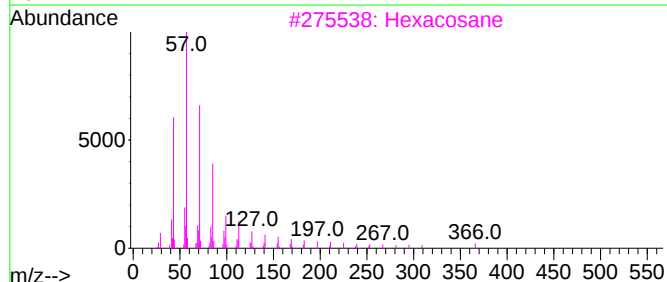
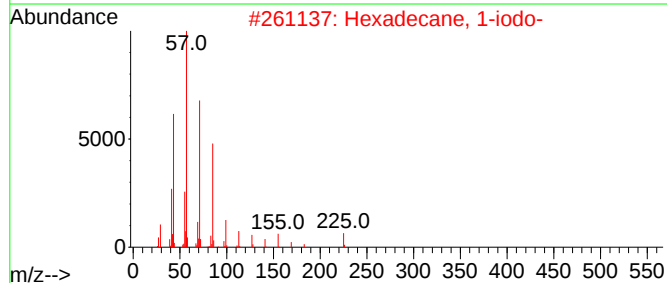
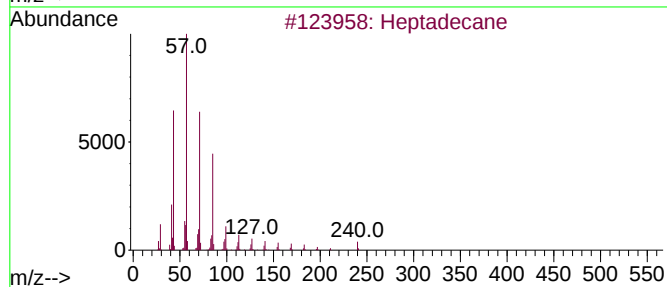
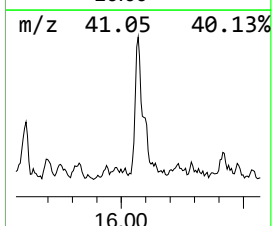
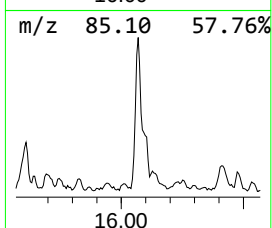
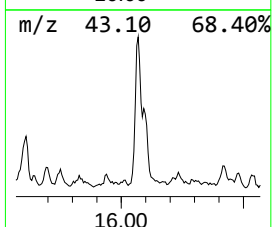
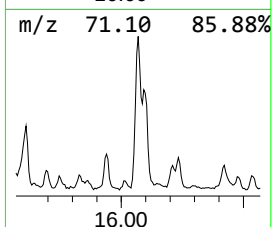
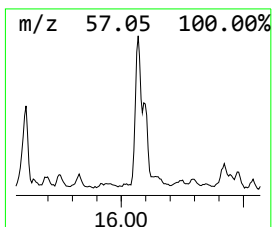
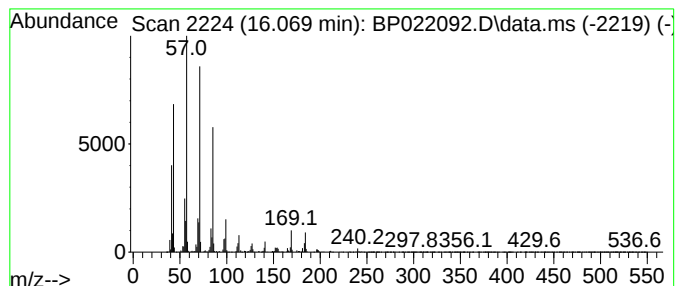
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Quant Title : SVOA CALIBRATION

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

Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.069	7.09 ng/ul	675323	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	80
2			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	80
3			Hexacosane	366	C26H54	000630-01-3	80
4			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	74
5			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
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Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

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ClientSampleId :
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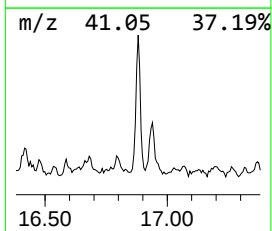
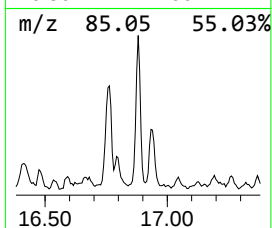
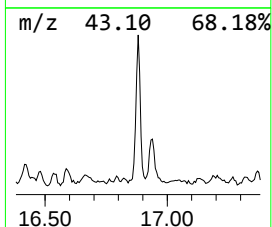
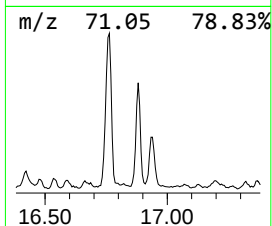
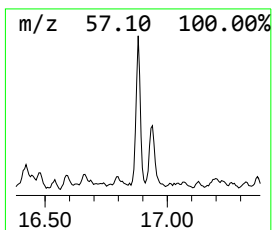
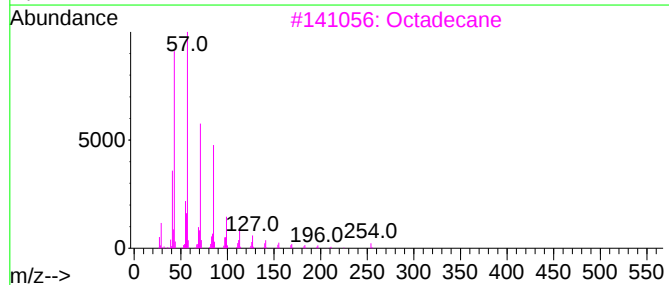
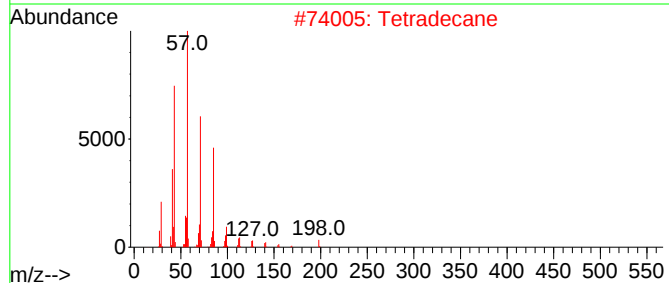
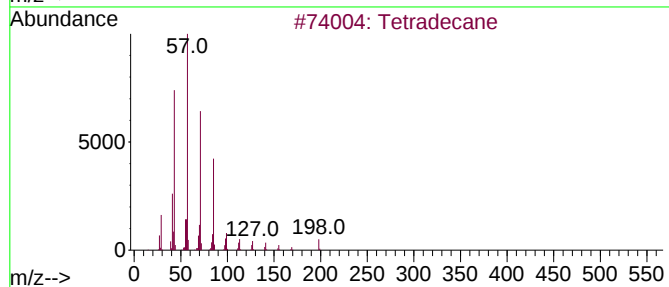
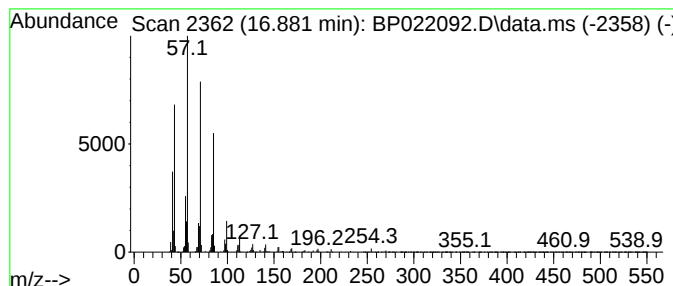
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Quant Title : SVOA CALIBRATION

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

Peak Number 52 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.881	5.52 ng/ul	525355	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	91
2			Tetradecane	198	C14H30	000629-59-4	90
3			Octadecane	254	C18H38	000593-45-3	89
4			Pentadecane	212	C15H32	000629-62-9	80
5			Heptadecane	240	C17H36	000629-78-7	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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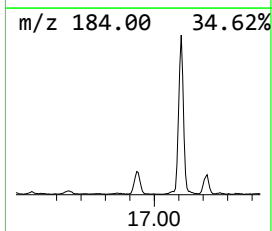
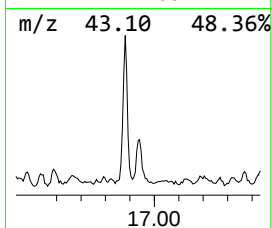
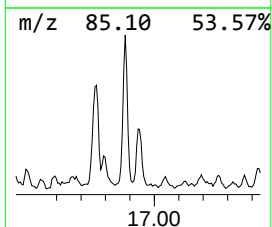
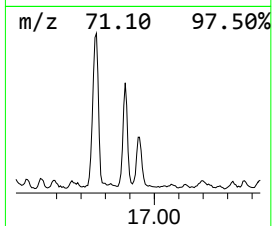
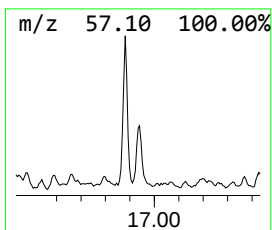
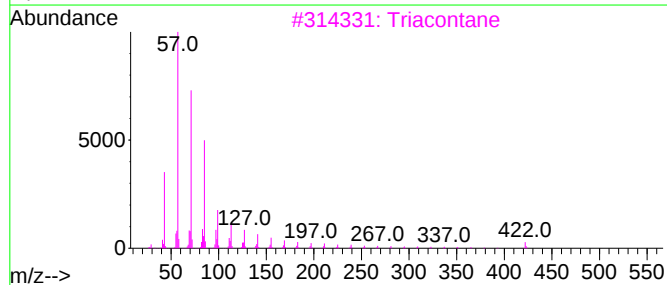
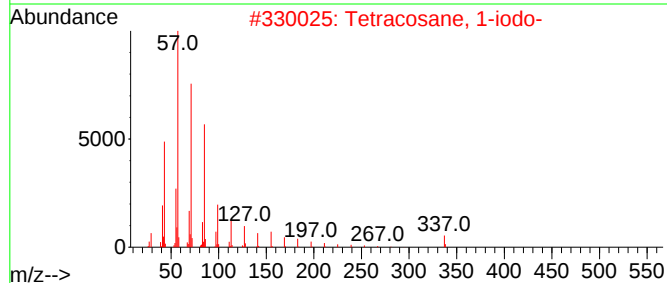
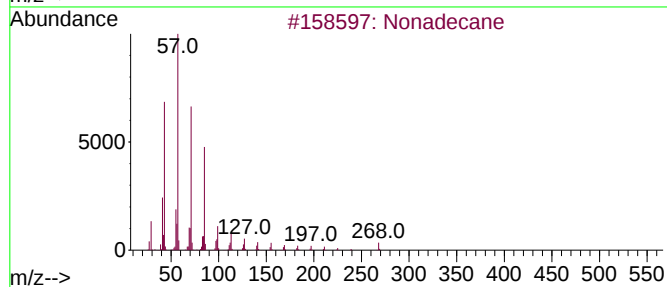
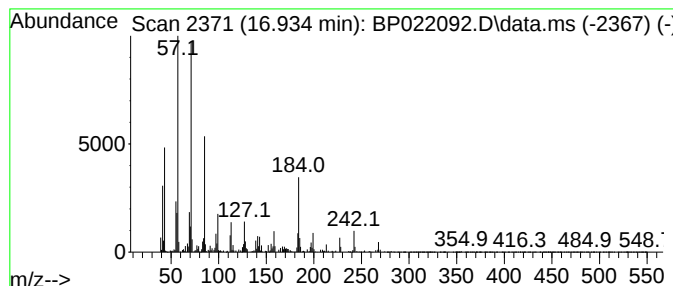
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Quant Title : SVOA CALIBRATION

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

Peak Number 53 (DEL) Alkane: Straight-Chai... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.934	4.28 ng/ul	407949	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	81
2			Tetracosane, 1-iodo-	464	C24H49I	1000406-32-0	72
3			Triacontane	422	C30H62	000638-68-6	72
4			2-Bromotetradecane	276	C14H29Br	074036-95-6	72
5			Docosane, 1-iodo-	436	C22H45I	1000406-31-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC63

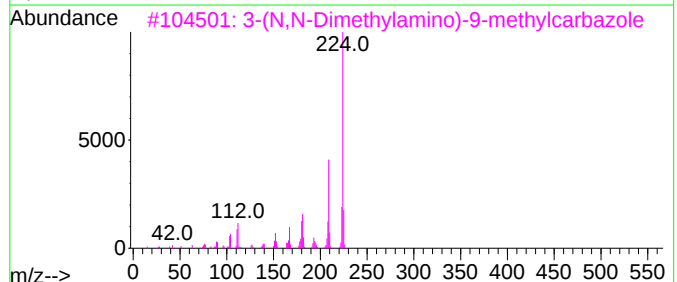
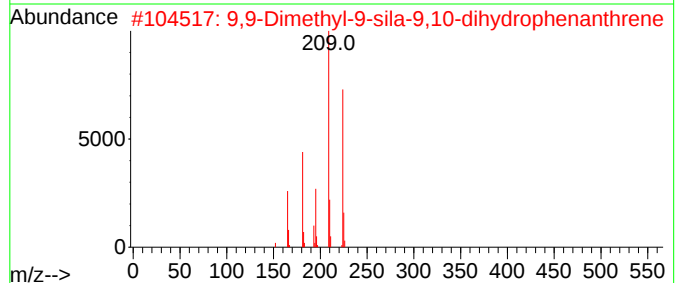
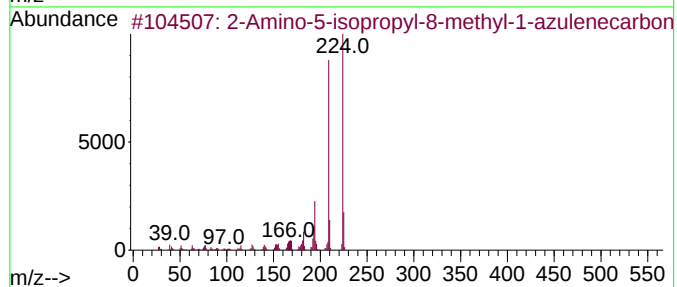
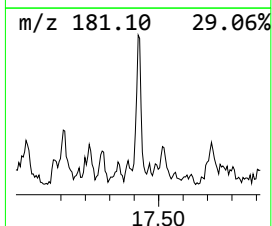
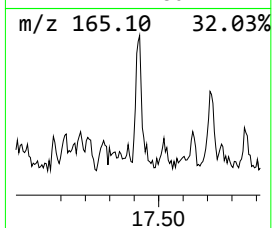
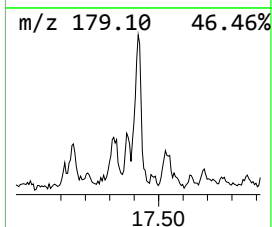
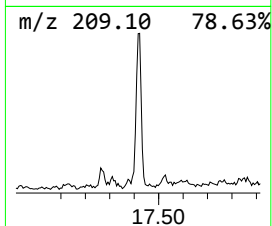
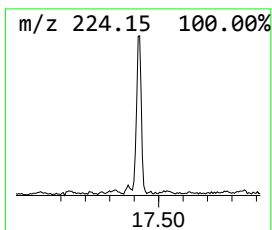
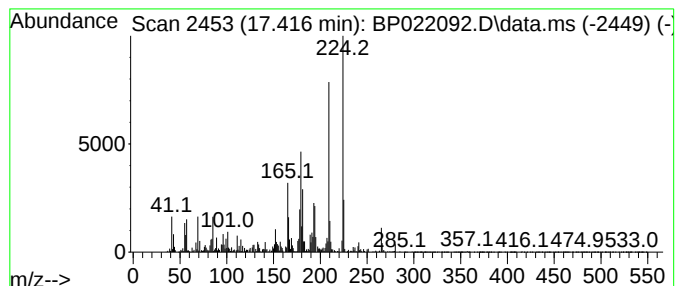
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Quant Title : SVOA CALIBRATION

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

Peak Number 54 2-Amino-5-isopropyl-8-methy... Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.416	3.41 ng/ul	324449	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Amino-5-isopropyl-8-methyl-1-a...	224	C15H16N2	093946-48-6	83
2			9,9-Dimethyl-9-sila-9,10-dihydro...	224	C15H16Si	187328-55-8	58
3			3-(N,N-Dimethylamino)-9-methylca...	224	C15H16N2	094127-15-8	38
4			1,2,3,4-Tetrahydronaphthalen-1,2...	224	C12H16O4	1000125-87-6	38
5			Benzoic acid, 2-(2-phenylethenyl)-	224	C15H12O2	060901-21-5	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC63

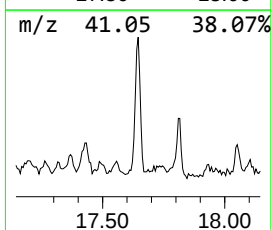
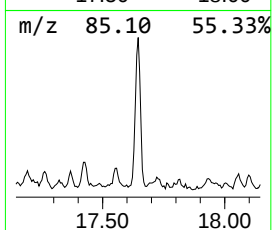
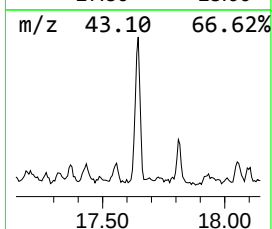
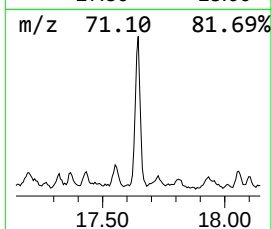
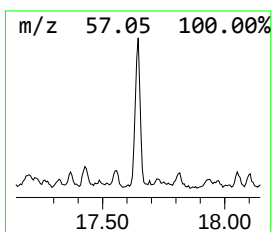
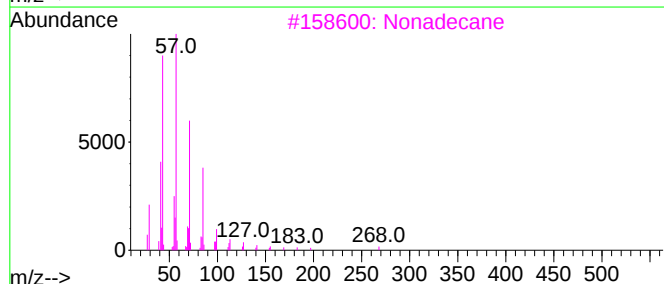
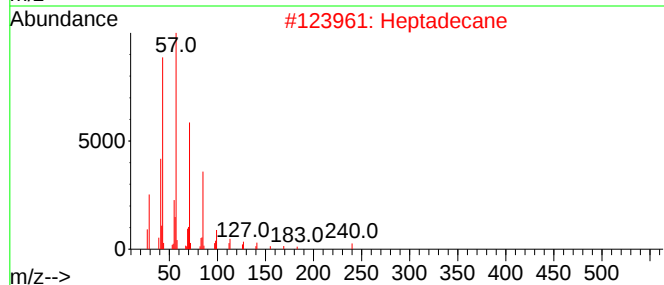
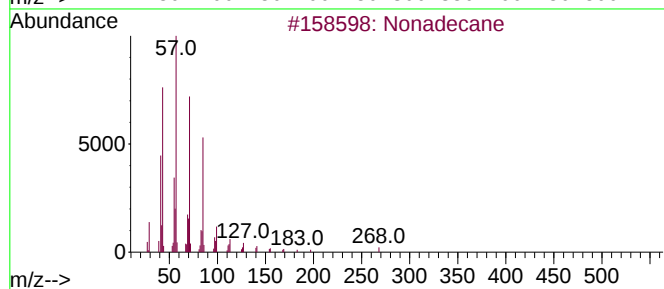
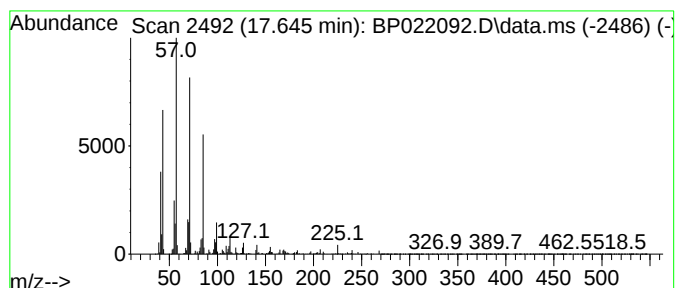
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Quant Title : SVOA CALIBRATION

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.
17.645	4.84 ng/ul	460772	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	95
2			Heptadecane	240	C17H36	000629-78-7	95
3			Nonadecane	268	C19H40	000629-92-5	90
4			Hexadecane	226	C16H34	000544-76-3	87
5			Dodecane	170	C12H26	000112-40-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC63

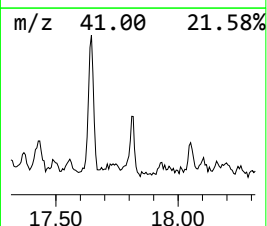
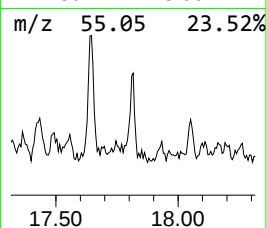
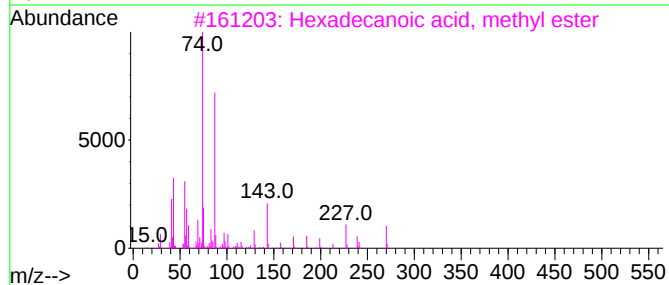
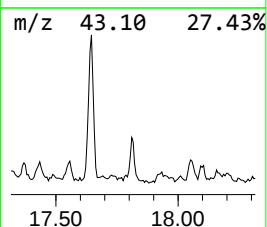
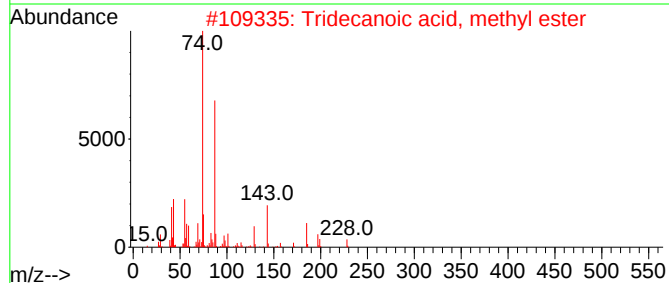
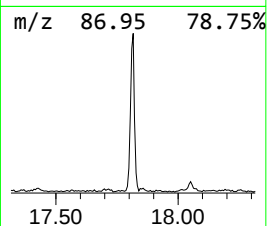
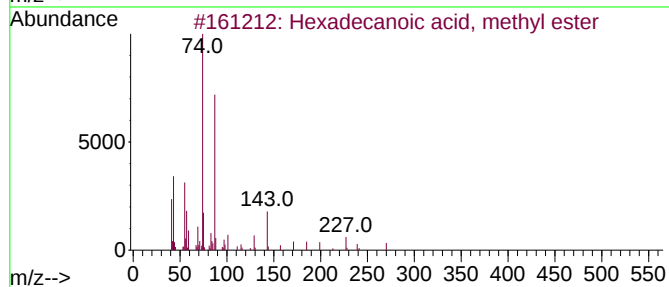
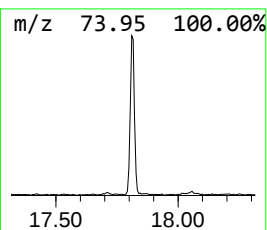
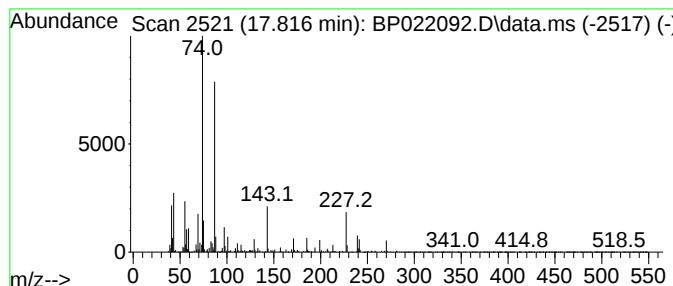
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Quant Title : SVOA CALIBRATION

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

Peak Number 56 Hexadecanoic acid, methyl e... Concentration Rank 60

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.816	2.06 ng/ul	195832	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	93
2			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	91
3			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91
4			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	87
5			Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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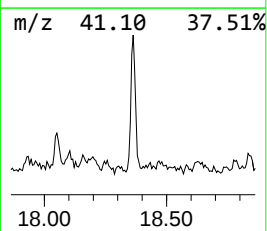
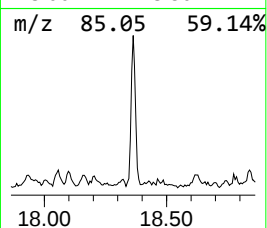
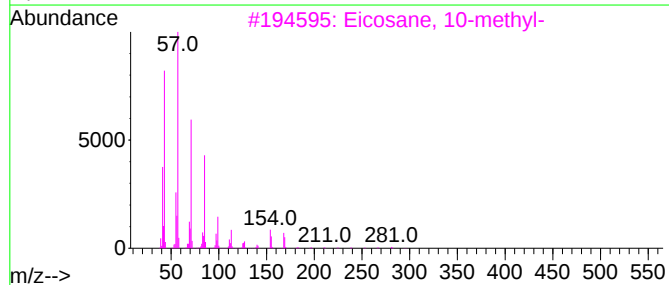
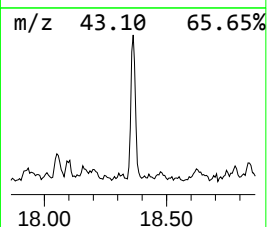
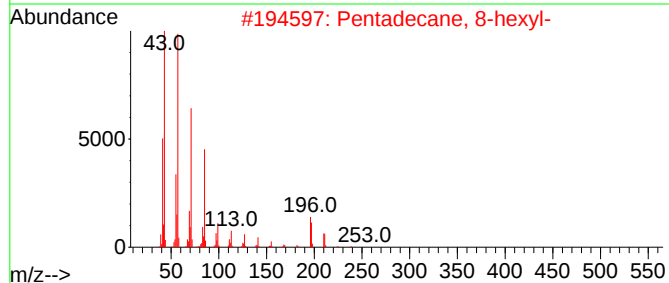
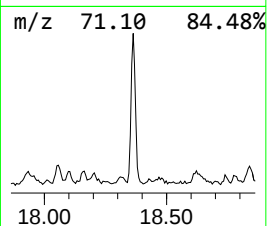
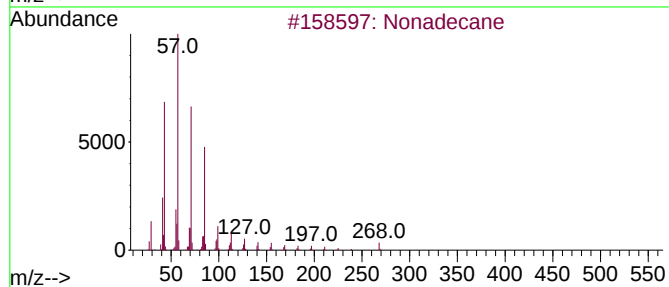
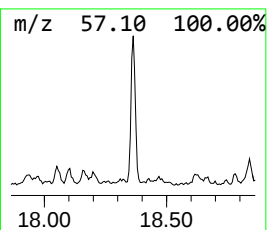
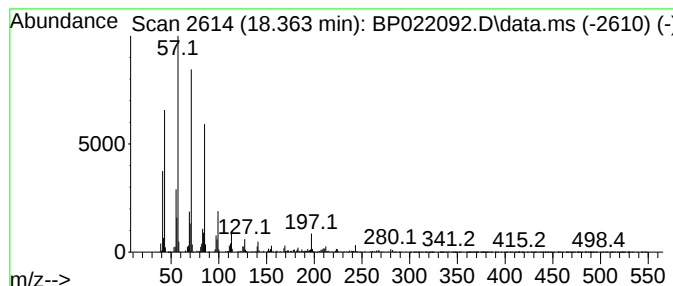
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Quant Title : SVOA CALIBRATION

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

Peak Number 57 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.363	3.81 ng/ul	363289	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	95
2			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
3			Eicosane, 10-methyl-	296	C21H44	054833-23-7	91
4			Octadecane	254	C18H38	000593-45-3	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC63

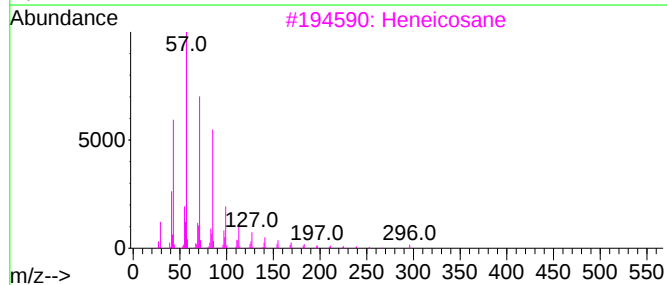
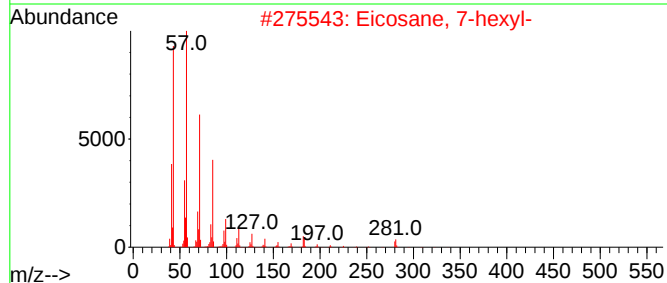
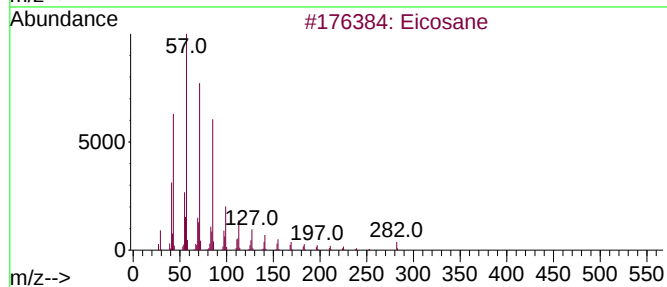
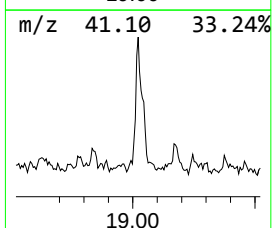
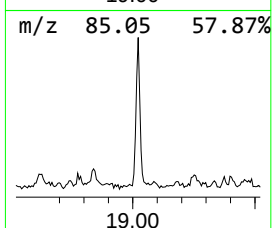
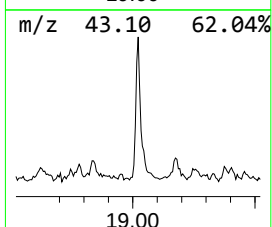
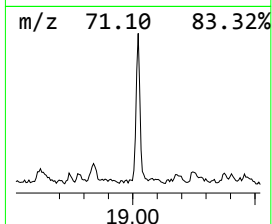
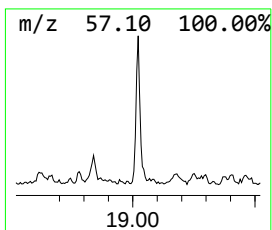
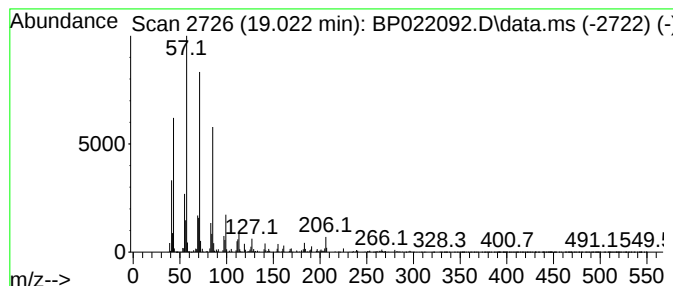
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Quant Title : SVOA CALIBRATION

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

Peak Number 58 (DEL) Alkane: Straight-Chai... Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.022	3.23 ng/ul	307231	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	96
2			Eicosane, 7-hexyl-	366	C26H54	055333-99-8	87
3			Heneicosane	296	C21H44	000629-94-7	87
4			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87
5			Eicosane	282	C20H42	000112-95-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC63

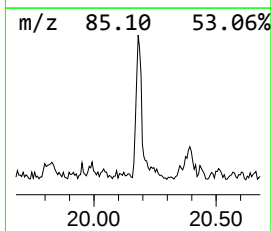
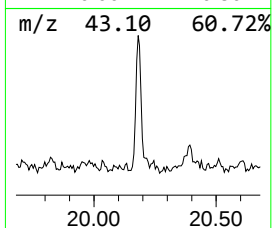
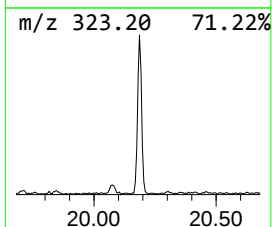
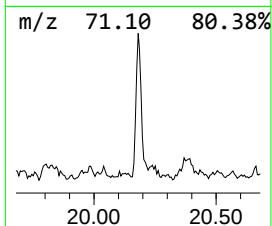
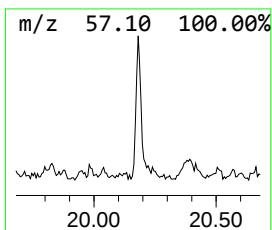
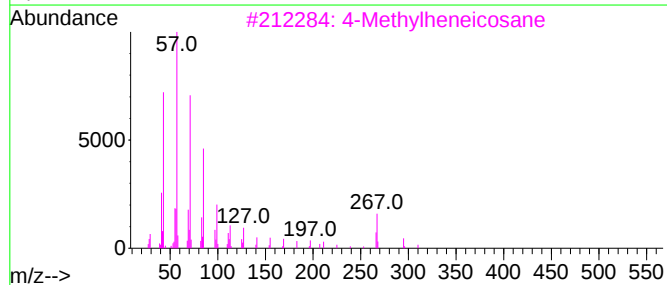
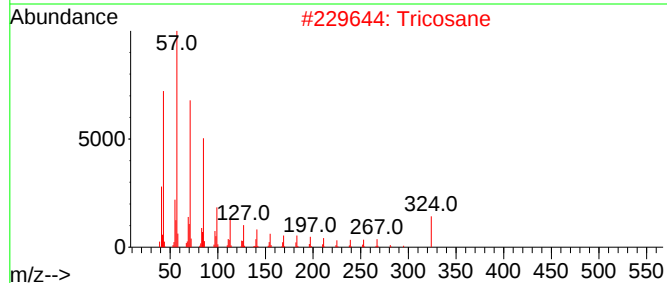
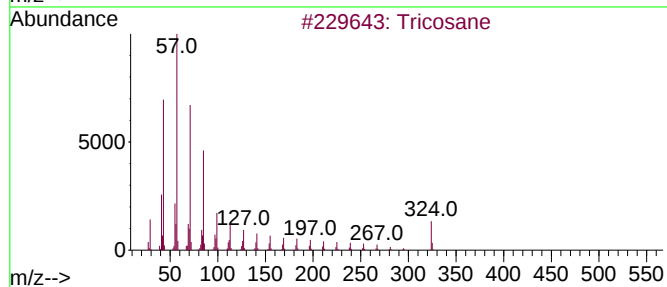
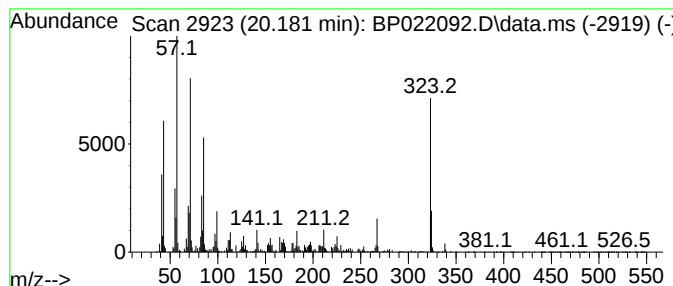
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Quant Title : SVOA CALIBRATION

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

Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.180	2.85 ng/ul	328608	Chrysene-d12	21.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tricosane	324	C23H48	000638-67-5	86
2			Tricosane	324	C23H48	000638-67-5	86
3			4-Methylheneicosane	310	C22H46	025117-29-7	55
4			5,5-Diethylheptadecane	296	C21H44	1010360-41-7	46
5			17,21-Dimethylheptatriacontane	549	C39H80	067979-79-7	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022092.D
Acq On : 28 Sep 2024 05:10
Operator : RC/JU
Sample : P3910-17 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC63

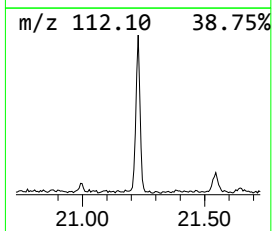
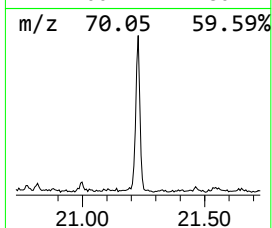
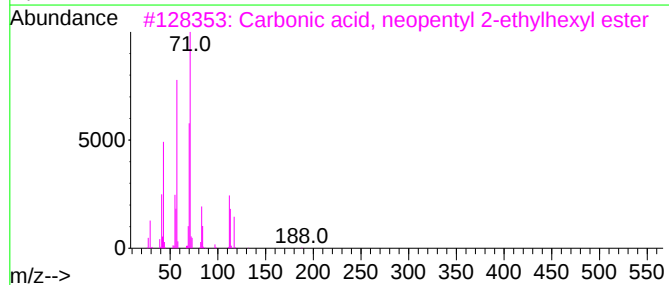
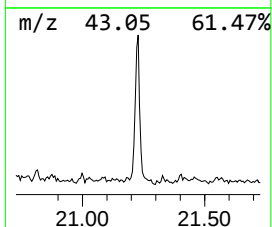
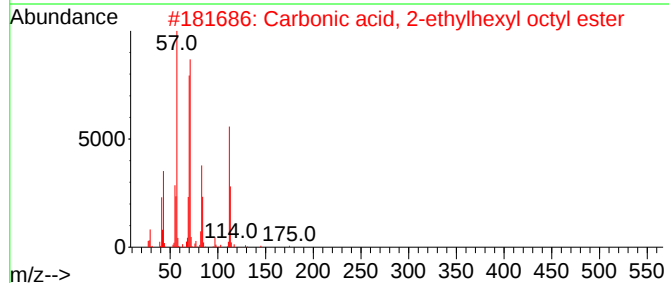
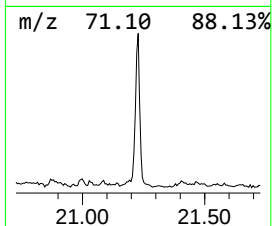
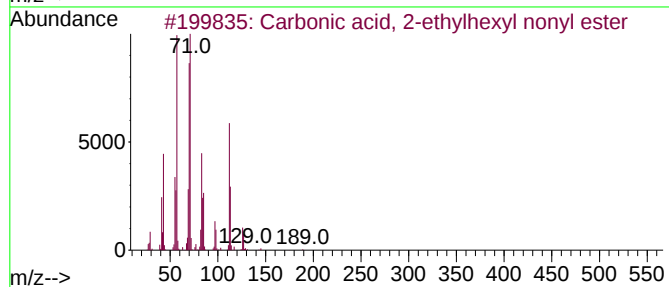
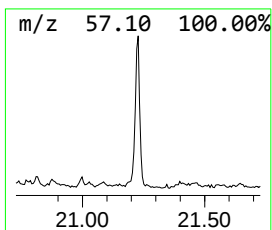
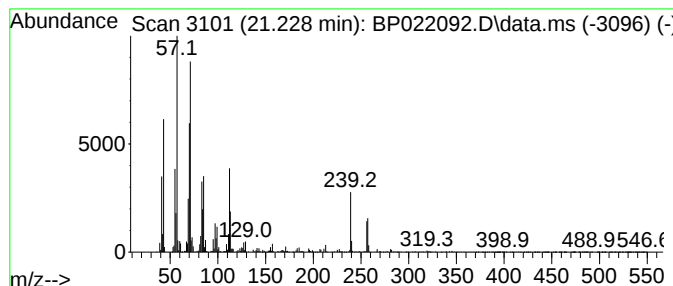
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 60 Carbonic acid, 2-ethylhexyl... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.227	4.34 ng/ul	500639	Chrysene-d12	21.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, 2-ethylhexyl nonyl...	300	C18H36O3	1000383-13-6	68
2			Carbonic acid, 2-ethylhexyl octy...	286	C17H34O3	1000383-13-5	64
3			Carbonic acid, neopentyl 2-ethyl...	244	C14H28O3	1000357-85-7	49
4			Carbonic acid, 2-ethylhexyl unde...	328	C20H40O3	1000383-13-8	46
5			Carbonic acid, decyl 2-ethylhexy...	314	C19H38O3	1000383-13-7	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
 Data File : BP022092.D
 Acq On : 28 Sep 2024 05:10
 Operator : RC/JU
 Sample : P3910-17 10X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DDC63

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
2-Pentanone, 4-...	4.870	3.0	ng/ul	147328	1	7.699	973668	20.0
(DEL) Alkane: S...	5.758	11.5	ng/ul	558495	1	7.699	973668	20.0
Hexylene glycol	6.058	8.1	ng/ul	394480	1	7.699	973668	20.0
unknown-01	6.217	2.8	ng/ul	136498	1	7.699	973668	20.0
(DEL) Alkane: S...	6.287	3.7	ng/ul	178818	1	7.699	973668	20.0
(DEL) Alkane: C...	6.369	4.1	ng/ul	197928	1	7.699	973668	20.0
(DEL) Alkane: S...	6.711	5.9	ng/ul	286505	1	7.699	973668	20.0
(DEL) Alkane: S...	6.769	5.9	ng/ul	285161	1	7.699	973668	20.0
Benzene, 1-ethy...	6.805	5.2	ng/ul	253837	1	7.699	973668	20.0
Benzene, 1-ethy...	7.099	4.1	ng/ul	199596	1	7.699	973668	20.0
2-Heptanone, 4,...	7.140	4.3	ng/ul	208352	1	7.699	973668	20.0
(DEL) Alkane: S...	7.334	36.8	ng/ul	1789440	1	7.699	973668	20.0
Tetradecane, 1-...	7.622	4.2	ng/ul	203172	1	7.699	973668	20.0
1-Hexanol, 2-et...	7.764	9.8	ng/ul	476504	1	7.699	973668	20.0
Benzene, 1,2,3-...	7.816	4.1	ng/ul	200038	1	7.699	973668	20.0
(DEL) Alkane: C...	7.987	3.8	ng/ul	184899	1	7.699	973668	20.0
Cyclohexanone, ...	8.122	15.2	ng/ul	739637	1	7.699	973668	20.0
(DEL) Alkane: S...	8.228	6.0	ng/ul	293152	1	7.699	973668	20.0
(DEL) Alkane: S...	8.281	2.9	ng/ul	139332	1	7.699	973668	20.0
Benzene, 1,3-di...	8.334	6.0	ng/ul	292611	1	7.699	973668	20.0
(DEL) Alkane: S...	8.458	5.9	ng/ul	285996	1	7.699	973668	20.0
Benzene, 1-meth...	8.493	4.6	ng/ul	222389	1	7.699	973668	20.0
Benzene, 2-ethy...	8.640	3.5	ng/ul	171386	1	7.699	973668	20.0
p-Cymene	8.687	6.1	ng/ul	295972	1	7.699	973668	20.0
(DEL) Alkane: S...	8.916	29.6	ng/ul	1441230	1	7.699	973668	20.0
unknown-02	9.034	6.9	ng/ul	335421	1	7.699	973668	20.0
2,4-Dipropyl-5-...	10.840	5.7	ng/ul	863178	2	10.463	3016090	20.0
(DEL) Alkane: S...	11.487	2.2	ng/ul	334046	2	10.463	3016090	20.0
Benzene, 1-(2-b...	11.552	4.0	ng/ul	608520	2	10.463	3016090	20.0
(DEL) Alkane: S...	11.887	5.1	ng/ul	766665	2	10.463	3016090	20.0
(DEL) Alkane: S...	12.852	2.2	ng/ul	171563	3	14.316	1575040	20.0
(DEL) Alkane: S...	13.134	8.0	ng/ul	625958	3	14.316	1575040	20.0
Naphthalene, 1,...	13.487	4.4	ng/ul	348986	3	14.316	1575040	20.0
Naphthalene, 2,...	13.622	3.7	ng/ul	287869	3	14.316	1575040	20.0
(DEL) Alkane: S...	13.793	3.5	ng/ul	271919	3	14.316	1575040	20.0
Naphthalene, 2,...	13.863	2.7	ng/ul	211442	3	14.316	1575040	20.0
unknown-03	13.946	2.9	ng/ul	228621	3	14.316	1575040	20.0
(DEL) Alkane: S...	14.228	11.3	ng/ul	888690	3	14.316	1575040	20.0
(DEL) Alkane: C...	14.399	3.1	ng/ul	246518	3	14.316	1575040	20.0
9-Methyl-S-octa...	14.457	9.7	ng/ul	762451	3	14.316	1575040	20.0
Heptacosane, 1-...	14.857	2.8	ng/ul	219635	3	14.316	1575040	20.0
2H-1-Benzopyran...	15.010	2.4	ng/ul	184741	3	14.316	1575040	20.0
Naphthalene, 1,...	15.081	11.0	ng/ul	869585	3	14.316	1575040	20.0
(DEL) Alkane: S...	15.187	7.0	ng/ul	553580	3	14.316	1575040	20.0
(DEL) Alkane: S...	15.610	3.8	ng/ul	296125	3	14.316	1575040	20.0
Benzophenone	15.698	3.9	ng/ul	304814	3	14.316	1575040	20.0
unknown-04	15.981	3.8	ng/ul	359979	4	17.110	1904550	20.0
(DEL) Alkane: S...	16.069	7.1	ng/ul	675323	4	17.110	1904550	20.0
(DEL) Alkane: S...	16.881	5.5	ng/ul	525355	4	17.110	1904550	20.0
(DEL) Alkane: S...	16.934	4.3	ng/ul	407949	4	17.110	1904550	20.0
2-Amino-5-isopr...	17.416	3.4	ng/ul	324449	4	17.110	1904550	20.0
(DEL) Alkane: S...	17.645	4.8	ng/ul	460772	4	17.110	1904550	20.0

Hexadecanoic ac...	17.816	2.1	ng/ul	195832	4	17.110	1904550	20.0
(DEL) Alkane: S...	18.363	3.8	ng/ul	363289	4	17.110	1904550	20.0
(DEL) Alkane: S...	19.022	3.2	ng/ul	307231	4	17.110	1904550	20.0
(DEL) Alkane: S...	20.180	2.9	ng/ul	328608	5	21.539	2306630	20.0
Carbonic acid, ...	21.227	4.3	ng/ul	500639	5	21.539	2306630	20.0

Instrument :

BNA_P

ClientSampleId :

DDC63