

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
 Data File : BP020268.D
 Acq On : 09 May 2024 13:42
 Operator : MA/JU
 Sample : P2369-16DL 30X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43P0DL

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-BP050124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020268.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.652	431	436	441	rVB4	12815	21213	0.81%	0.125%
2	5.899	470	478	484	rBV9	10080	25269	0.97%	0.149%
3	6.011	491	497	502	rVB8	7959	17416	0.67%	0.102%
4	6.175	521	525	529	rBV	19753	25579	0.98%	0.150%
5	6.275	535	542	554	rVB3	23380	63809	2.44%	0.375%
6	6.505	573	581	586	rBV5	13236	32178	1.23%	0.189%
7	6.563	586	591	594	rBV5	10559	17244	0.66%	0.101%
8	6.599	594	597	602	rVB2	24682	33770	1.29%	0.199%
9	6.658	602	607	611	rBV2	21720	29871	1.14%	0.176%
10	6.705	611	615	618	rBV5	12009	17659	0.67%	0.104%
11	6.763	618	625	630	rBV3	41914	81741	3.12%	0.481%
12	6.846	634	639	643	rBV2	17933	27120	1.04%	0.159%
13	6.922	648	652	658	rVB4	16227	30812	1.18%	0.181%
14	6.999	658	665	669	rBV6	19032	53555	2.05%	0.315%
15	7.087	677	680	687	rVB3	22344	33917	1.30%	0.199%
16	7.163	688	693	698	rBV5	19275	35214	1.35%	0.207%
17	7.222	698	703	707	rBV	173326	243050	9.28%	1.429%
18	7.316	715	719	724	rVB3	13805	22364	0.85%	0.132%
19	7.399	729	733	735	rBV2	15570	24475	0.93%	0.144%
20	7.510	742	752	756	rBV5	47779	136406	5.21%	0.802%
21	7.563	756	761	765	rVV	142353	219039	8.37%	1.288%
22	7.616	765	770	782	rVB	323864	626050	23.92%	3.681%
23	7.722	782	788	792	rBV2	44980	95925	3.66%	0.564%
24	7.775	793	797	801	rVV4	60605	91331	3.49%	0.537%
25	7.840	805	808	811	rVV	37777	60889	2.33%	0.358%
26	7.875	811	814	820	rVB4	47625	71226	2.72%	0.419%
27	7.934	821	824	831	rVB8	23298	47346	1.81%	0.278%
28	8.022	835	839	846	rVB5	26993	52980	2.02%	0.312%
29	8.110	846	854	858	rBV3	91765	208958	7.98%	1.229%
30	8.169	861	864	869	rVB	77802	120246	4.59%	0.707%
31	8.240	869	876	881	rBV3	129736	241679	9.23%	1.421%
32	8.352	890	895	899	rBV	90891	145012	5.54%	0.853%
33	8.410	900	905	921	rVB7	81829	295382	11.28%	1.737%
34	8.546	923	928	932	rBV2	40846	80316	3.07%	0.472%
35	8.599	933	937	942	rVB2	44140	68905	2.63%	0.405%
36	8.657	942	947	949	rBV4	43991	74080	2.83%	0.436%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

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Stop Thrs : 0

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
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37	8.693	949	953	962	rVB5	66344	141276	5.40%	0.831%
38	8.810	963	973	982	rVB	561217	997659	38.11%	5.866%
39	8.893	982	987	989	rBV3	52178	90184	3.45%	0.530%
40	8.940	989	995	1002	rVB8	70702	172847	6.60%	1.016%
41	9.034	1006	1011	1019	rVV6	51020	116562	4.45%	0.685%
42	9.216	1038	1042	1046	rBV2	58741	101215	3.87%	0.595%
43	9.293	1049	1055	1059	rVV2	43354	101688	3.88%	0.598%
44	9.457	1077	1083	1087	rBV5	38352	87656	3.35%	0.515%
45	9.504	1087	1091	1096	rVV4	32623	84235	3.22%	0.495%
46	9.551	1096	1099	1103	rVV3	46829	78505	3.00%	0.462%
47	9.593	1103	1106	1111	rVV2	36478	72111	2.75%	0.424%
48	9.651	1114	1116	1121	rVB4	35440	47052	1.80%	0.277%
49	9.722	1123	1128	1132	rVV3	33882	60665	2.32%	0.357%
50	9.793	1135	1140	1146	rVV2	72623	150810	5.76%	0.887%
51	9.846	1147	1149	1155	rVB3	22265	32451	1.24%	0.191%
52	9.904	1155	1159	1163	rBV	26184	34756	1.33%	0.204%
53	9.963	1163	1169	1175	rVB5	28565	67900	2.59%	0.399%
54	10.093	1183	1191	1197	rBV8	20839	48805	1.86%	0.287%
55	10.151	1197	1201	1210	rVV8	10512	18749	0.72%	0.110%
56	10.251	1212	1218	1229	rVV8	15869	57478	2.20%	0.338%
57	10.351	1229	1235	1237	rVV	137602	214244	8.18%	1.260%
58	10.387	1237	1241	1255	rVV	485230	899073	34.35%	5.287%
59	10.499	1255	1260	1263	rVV4	33965	65192	2.49%	0.383%
60	10.534	1263	1266	1273	rVV	32735	47799	1.83%	0.281%
61	10.669	1285	1289	1297	rVB7	9330	22316	0.85%	0.131%
62	10.757	1297	1304	1310	rBV8	12217	27853	1.06%	0.164%
63	11.404	1410	1414	1418	rBV3	11621	17582	0.67%	0.103%
64	11.816	1479	1484	1489	rVV3	16270	26963	1.03%	0.159%
65	12.257	1553	1559	1573	rBV	29630	67622	2.58%	0.398%
66	12.504	1596	1601	1606	rVB3	10306	19185	0.73%	0.113%
67	12.775	1640	1647	1654	rBV3	6398	18877	0.72%	0.111%
68	13.104	1698	1703	1709	rBV8	12156	24373	0.93%	0.143%
69	13.340	1736	1743	1748	rVB4	17769	37786	1.44%	0.222%
70	13.451	1758	1762	1769	rVB7	13378	25260	0.96%	0.149%
71	13.640	1790	1794	1804	rVB7	13407	32314	1.23%	0.190%
72	13.728	1804	1809	1819	rBV2	27765	72982	2.79%	0.429%
73	13.916	1836	1841	1842	rBV3	15936	22701	0.87%	0.133%
74	13.939	1842	1845	1849	rVV5	19895	27781	1.06%	0.163%
75	13.992	1849	1854	1859	rVV2	23733	46674	1.78%	0.274%
76	14.051	1859	1864	1869	rVV	33779	55322	2.11%	0.325%

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77	14.128	1872	1877	1886	rVB5	17555	41147	1.57%	0.242%
78	14.292	1899	1905	1913	rVV	712025	1211511	46.28%	7.124%
79	14.357	1913	1916	1922	rVB3	35115	50257	1.92%	0.296%
80	14.539	1938	1947	1951	rBV	91343	153943	5.88%	0.905%
81	14.892	2001	2007	2013	rVV5	21399	37907	1.45%	0.223%
82	14.963	2013	2019	2029	rVV6	38757	82427	3.15%	0.485%
83	15.157	2048	2052	2060	rVB3	17833	29915	1.14%	0.176%
84	15.334	2074	2082	2088	rBV2	39114	70373	2.69%	0.414%
85	15.416	2093	2096	2107	rVB8	17300	34697	1.33%	0.204%
86	15.598	2123	2127	2132	rBV7	12146	23011	0.88%	0.135%
87	16.034	2193	2201	2204	rBV8	9811	24221	0.93%	0.142%
88	16.769	2320	2326	2344	rBV	1360020	2617755	100.00%	15.393%
89	16.916	2349	2351	2358	rVB7	9792	18664	0.71%	0.110%
90	17.133	2381	2388	2397	rBV	888930	1428595	54.57%	8.400%
91	17.251	2403	2408	2413	rBV2	31920	50530	1.93%	0.297%
92	19.069	2713	2717	2720	rBV	14773	19231	0.73%	0.113%
93	19.657	2812	2817	2819	rVV2	39859	69557	2.66%	0.409%
94	19.680	2819	2821	2832	rVB3	58620	82532	3.15%	0.485%
95	20.251	2914	2918	2924	rBV	75410	106585	4.07%	0.627%
96	20.774	3003	3007	3012	rBV	98975	135124	5.16%	0.795%
97	21.316	3095	3099	3105	rVB2	84694	123989	4.74%	0.729%
98	21.633	3147	3153	3165	rVB2	668887	1290162	49.29%	7.586%
99	21.898	3195	3198	3204	rVB2	71205	100701	3.85%	0.592%
100	24.998	3716	3725	3744	rVB	470167	1397296	53.38%	8.216%

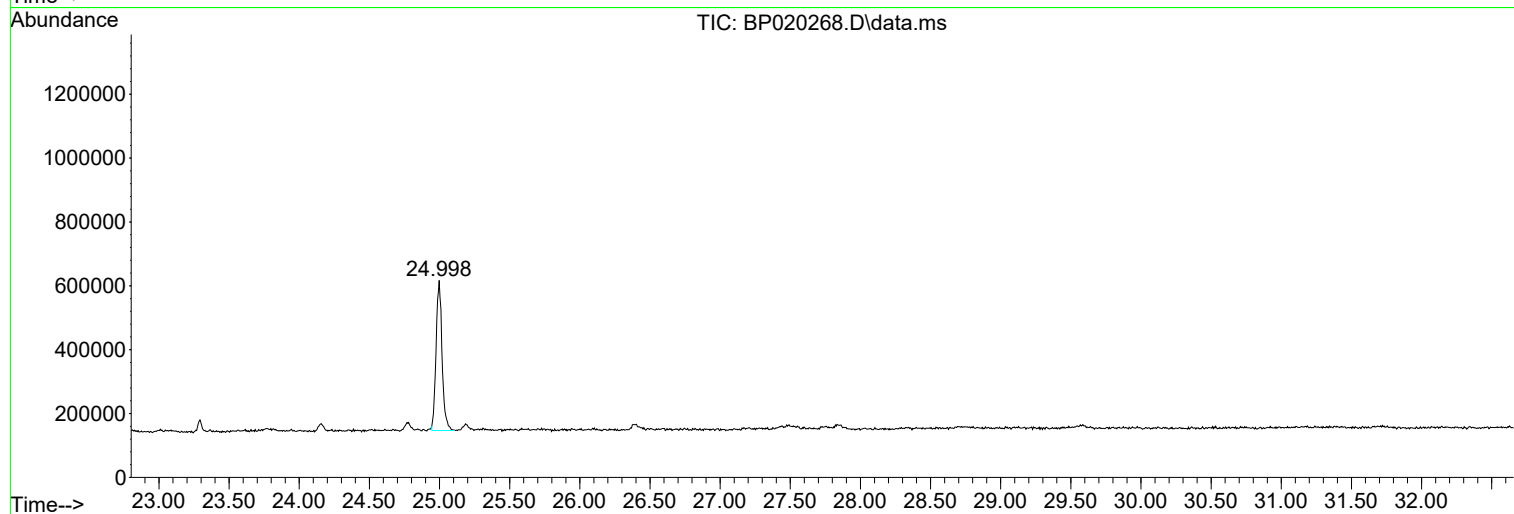
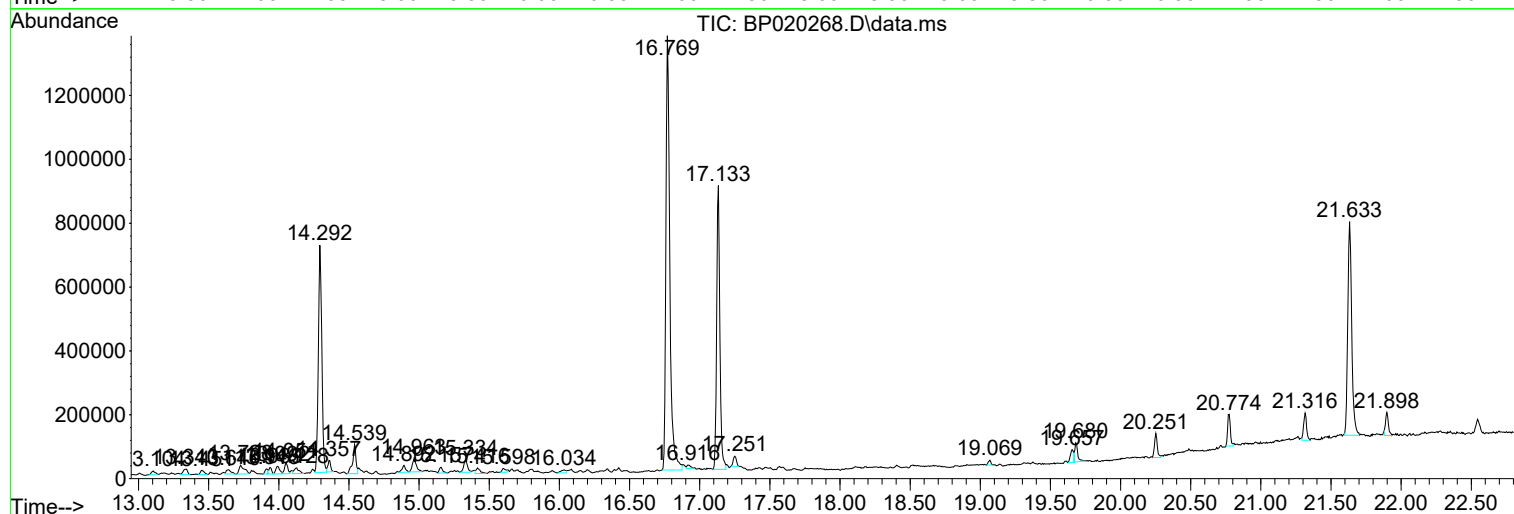
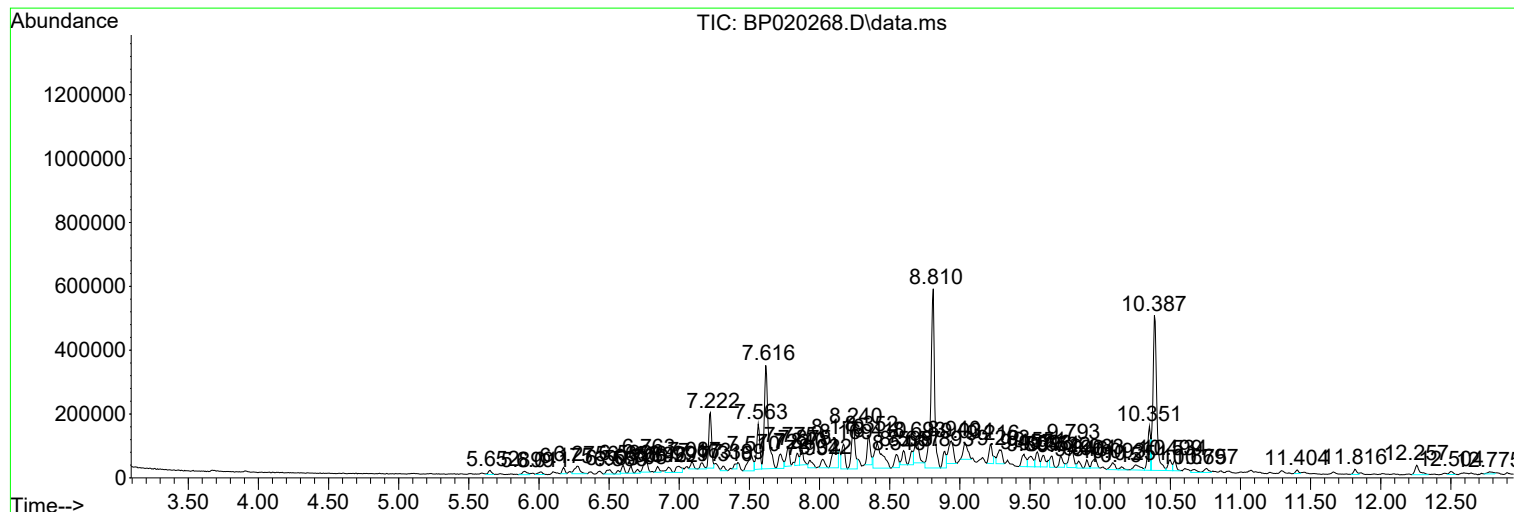
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ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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



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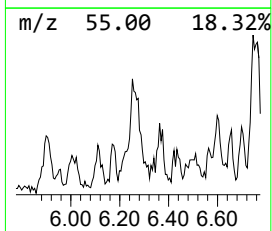
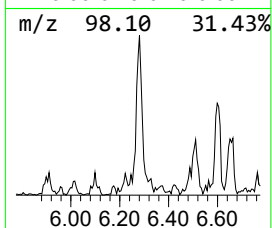
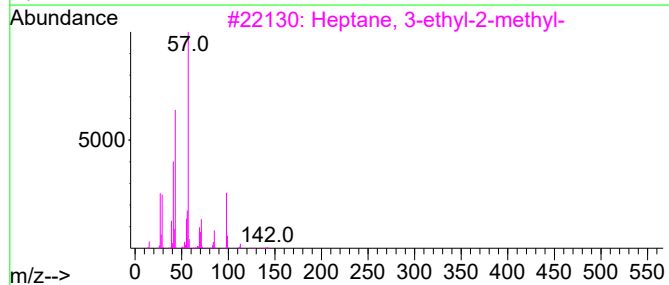
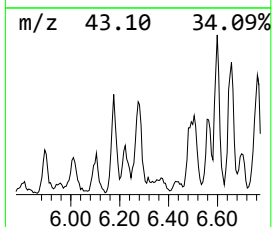
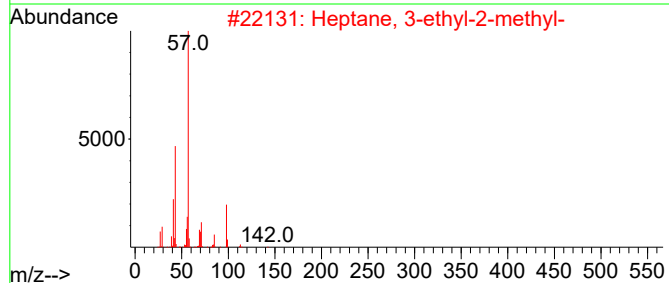
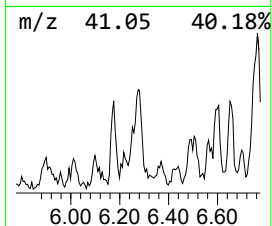
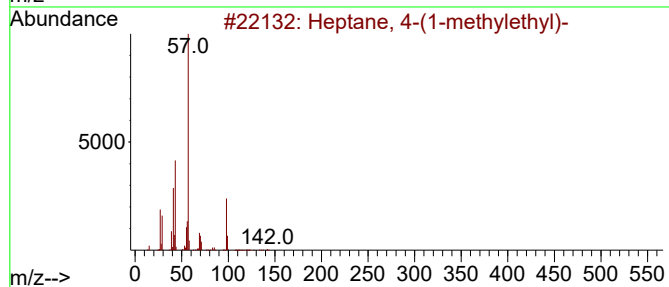
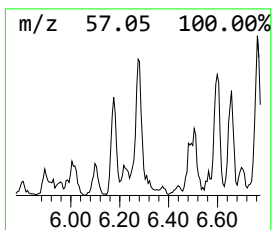
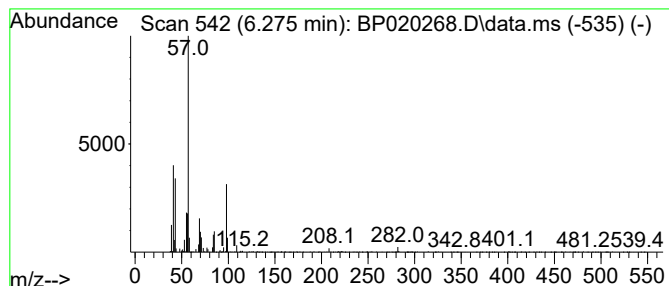
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.275	2.04 ng/ul	63809	1,4-Dichlorobenzene-d4	7.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 4-(1-methylethyl)-	142	C10H22	052896-87-4	52
2			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	50
3			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	50
4			Decane, 2,5-dimethyl-	170	C12H26	017312-50-4	50
5			Octane, 4,5-dipropyl-	198	C14H30	020905-05-9	45



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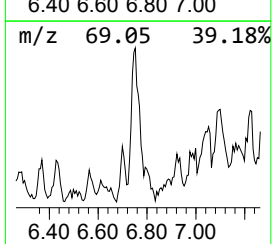
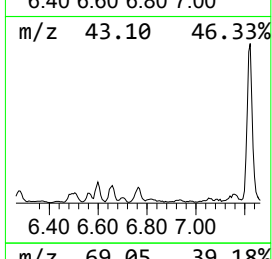
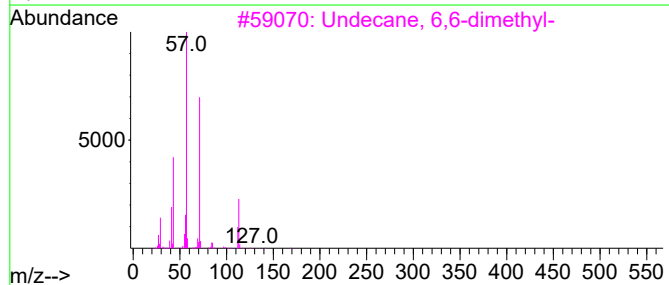
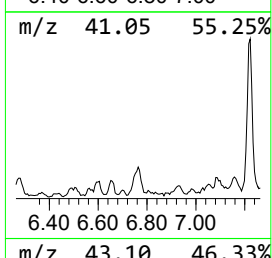
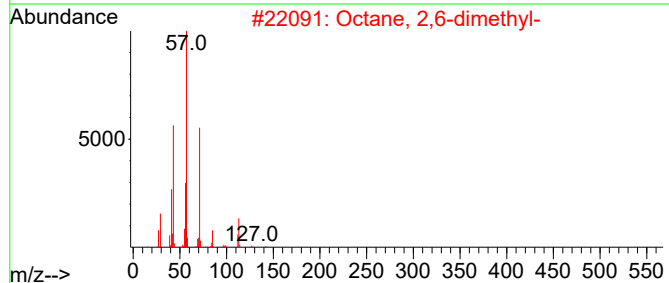
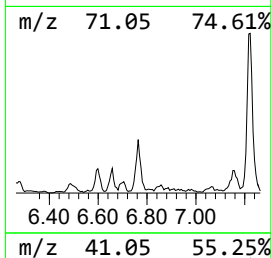
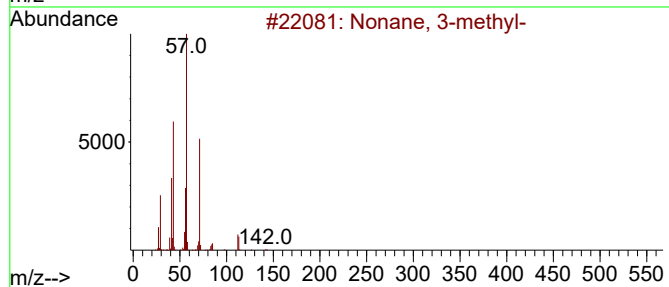
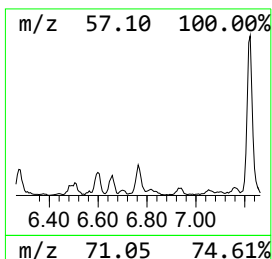
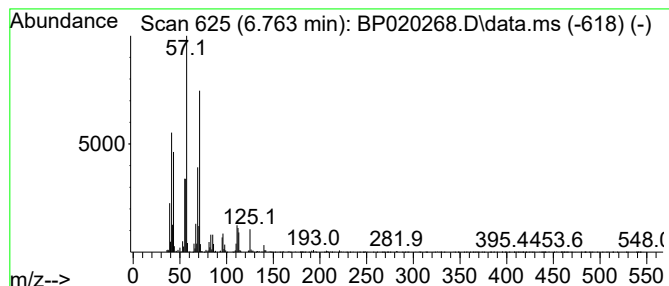
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD			R.T.
6.763	2.61 ng/ul	81741	1,4-Dichlorobenzene-d4			7.616
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 3-methyl-		142	C10H22	005911-04-6	72
2	Octane, 2,6-dimethyl-		142	C10H22	002051-30-1	68
3	Undecane, 6,6-dimethyl-		184	C13H28	017312-76-4	53
4	Butane, 2,2-dimethyl-		86	C6H14	000075-83-2	53
5	Threitol, 2-O-octyl-		234	C12H26O4	163776-14-5	53



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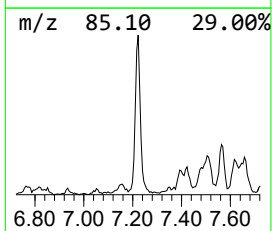
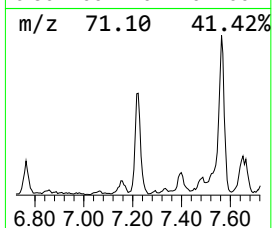
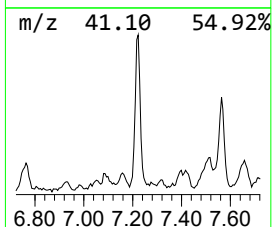
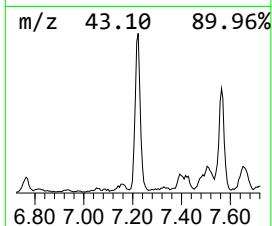
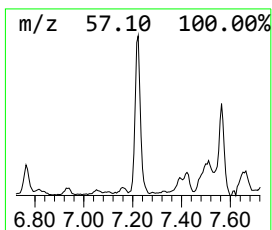
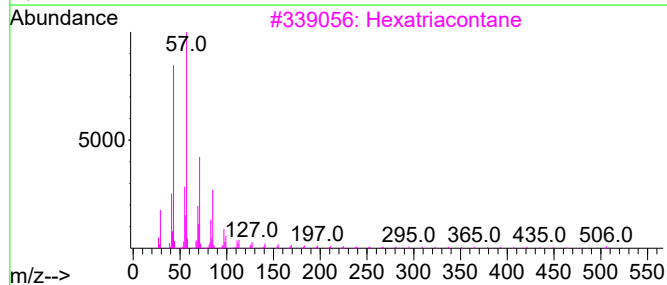
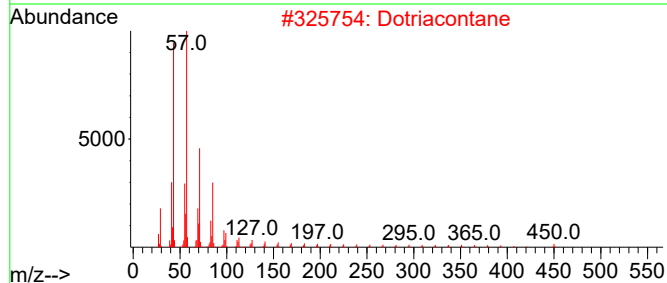
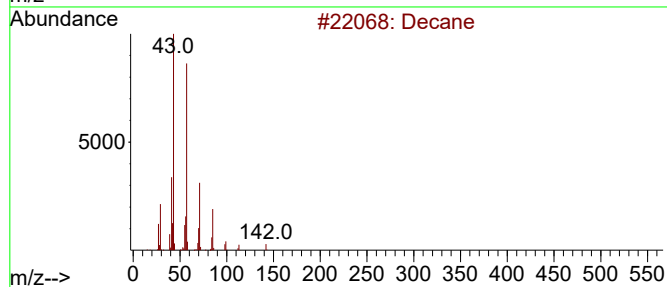
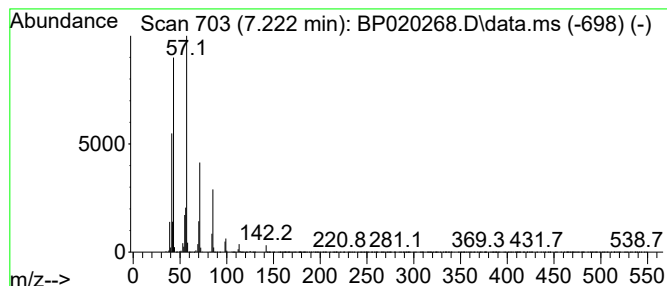
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.222	7.76 ng/ul	243050	1,4-Dichlorobenzene-d4	7.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	91
2			Dotriacontane	451	C32H66	000544-85-4	83
3			Hexatriacontane	507	C36H74	000630-06-8	83
4			Decane	142	C10H22	000124-18-5	78
5			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
Operator : MA/JU
Sample : P2369-16DL 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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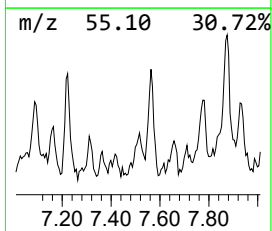
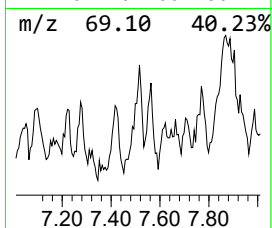
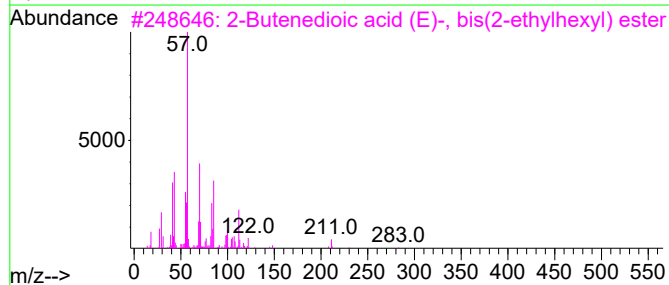
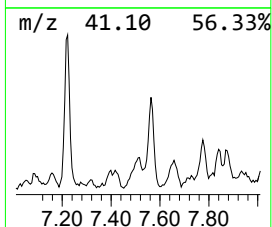
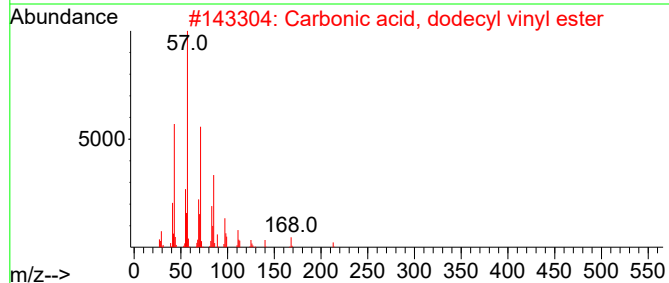
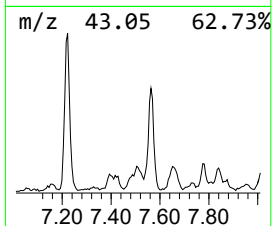
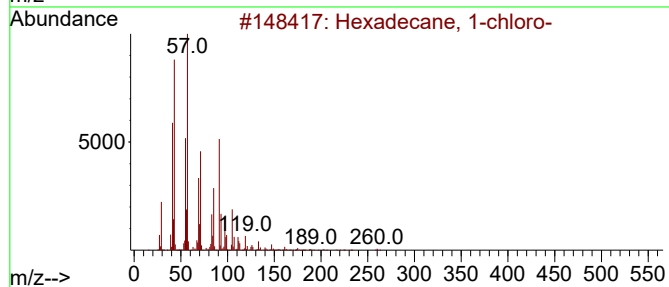
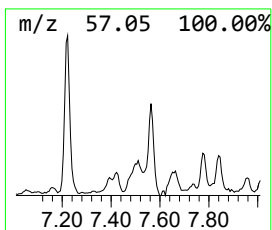
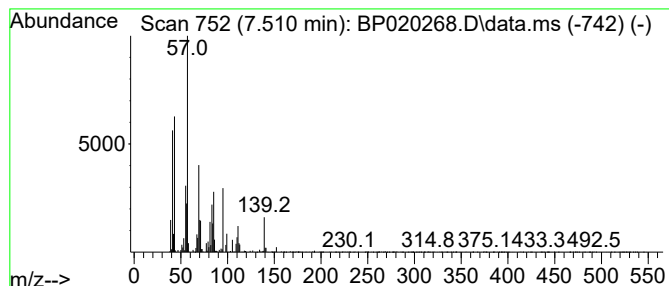
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Hexadecane, 1-chloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.510	4.36 ng/ul	136406	1,4-Dichlorobenzene-d4	7.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	59
2			Carbonic acid, dodecyl vinyl ester	256	C15H28O3	1000382-54-8	53
3			2-Butenedioic acid (E)-, bis(2-e...	340	C20H36O4	000141-02-6	50
4			Cyclobutylcarboxamide, N-allyl-	139	C8H13NO	1000340-36-9	49
5			6-Undecen-3-one, 5-butyl-2,2-dim...	252	C17H32O	055976-05-1	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
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Instrument :
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ClientSampleId :
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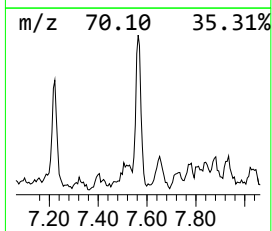
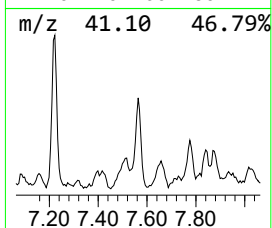
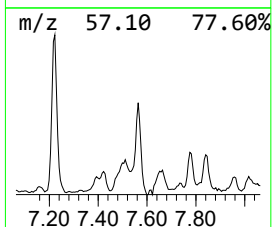
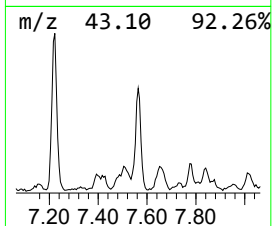
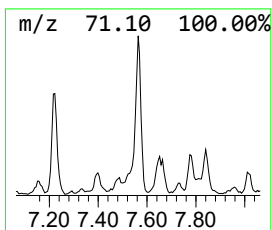
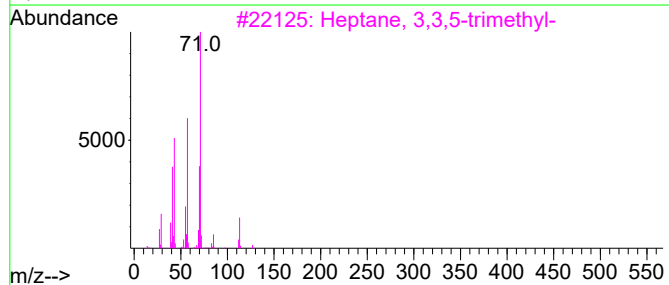
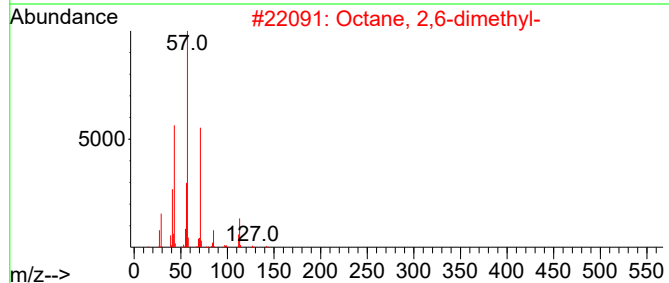
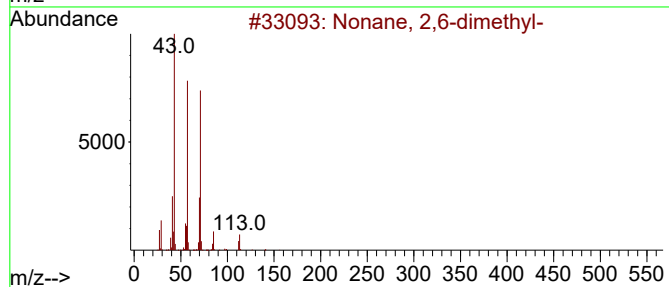
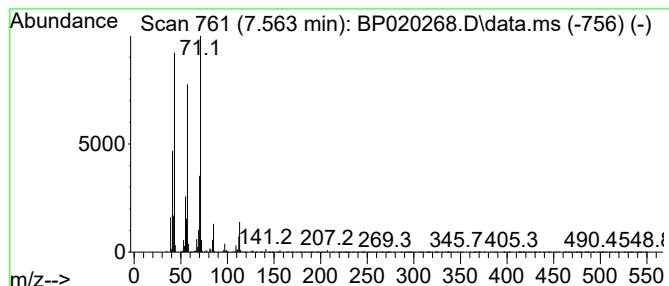
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.563	7.00 ng/ul	219039	1,4-Dichlorobenzene-d4	7.616

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	91
2	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	86
3	Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	78
4	Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	72
5	Sulfurous acid, dodecyl 2-pentyl...	320	C17H36O3S	1000309-16-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
Operator : MA/JU
Sample : P2369-16DL 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
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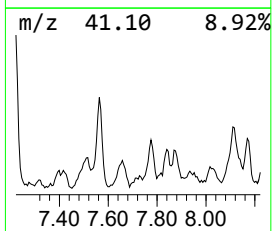
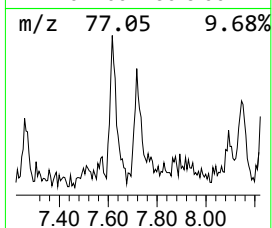
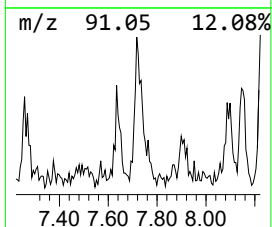
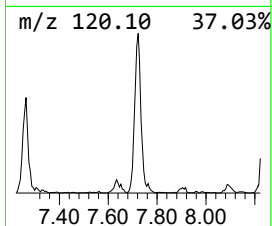
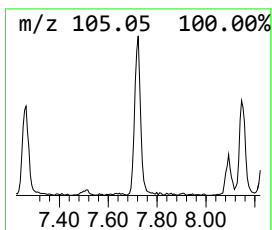
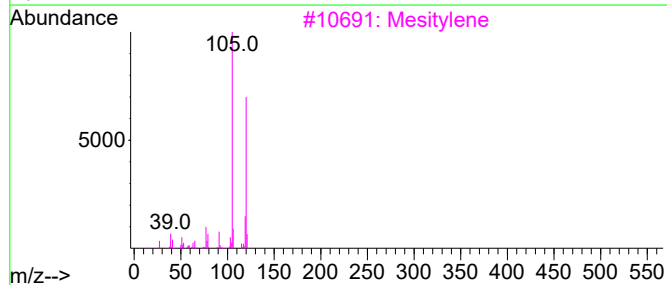
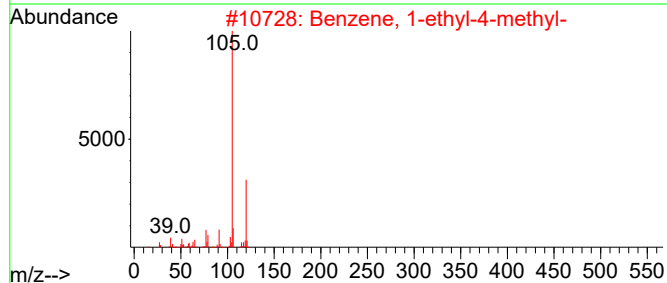
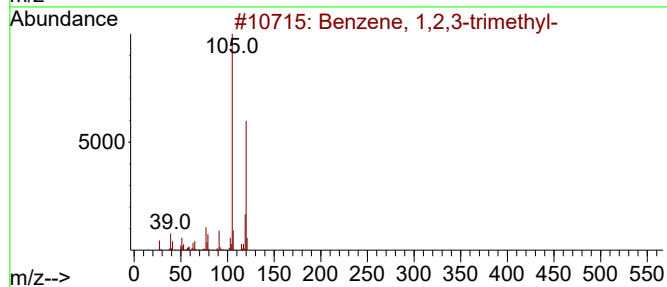
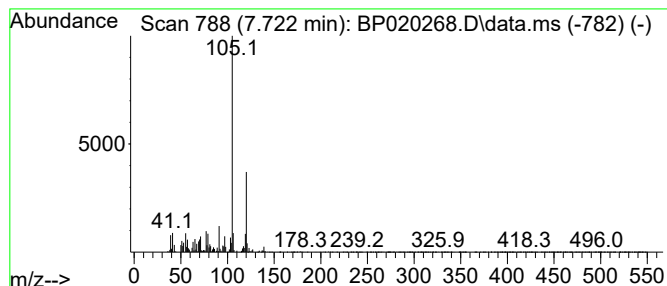
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.722	3.06 ng/ul	95925	1,4-Dichlorobenzene-d4	7.616

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
Operator : MA/JU
Sample : P2369-16DL 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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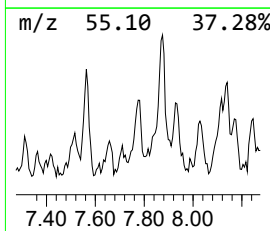
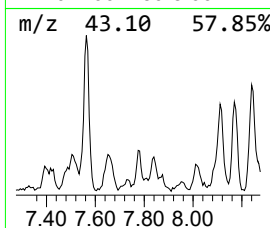
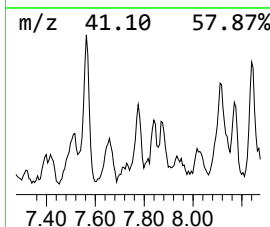
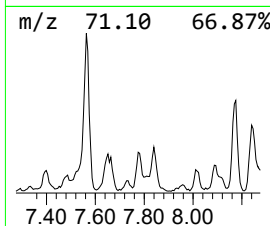
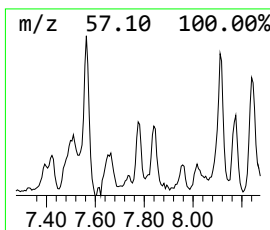
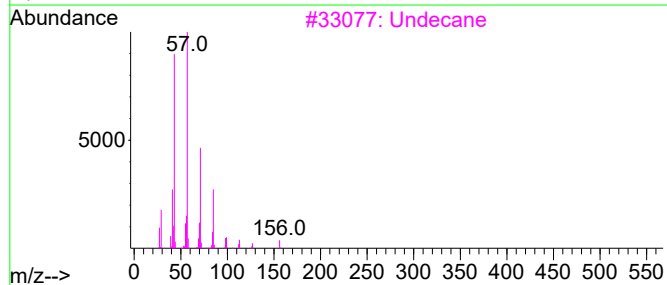
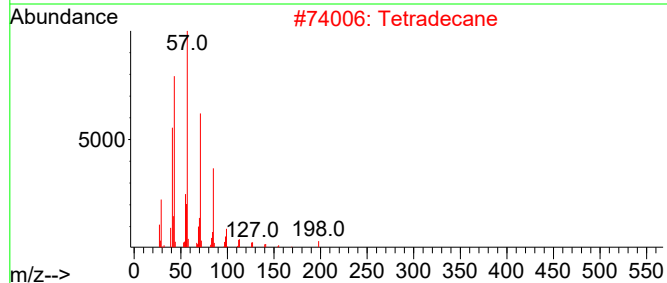
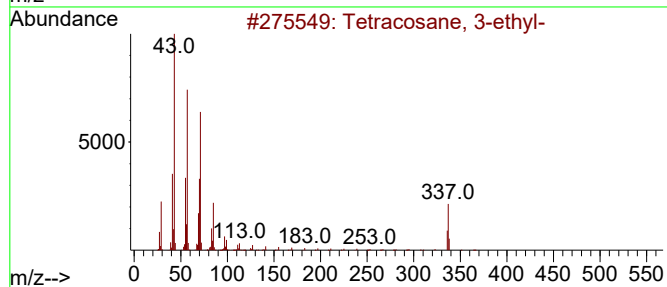
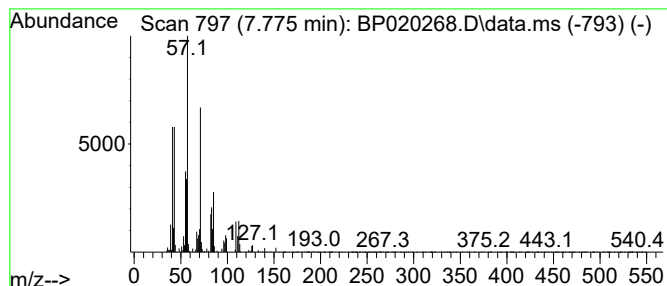
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.775	2.92 ng/ul	91331	1,4-Dichlorobenzene-d4	7.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane, 3-ethyl-	366	C26H54	055282-17-2	58
2			Tetradecane	198	C14H30	000629-59-4	52
3			Undecane	156	C11H24	001120-21-4	50
4			Bacchotricuneatin c	342	C20H22O5	066563-30-2	49
5			Nonadecane, 9-methyl-	282	C20H42	013287-24-6	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
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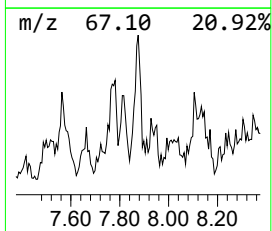
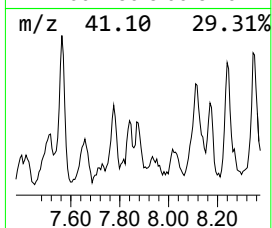
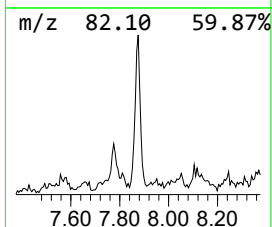
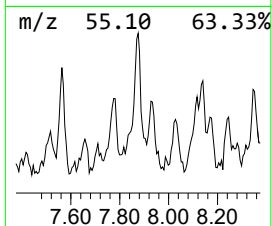
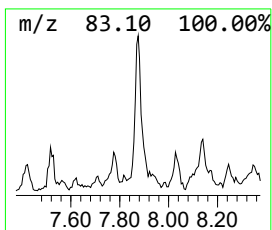
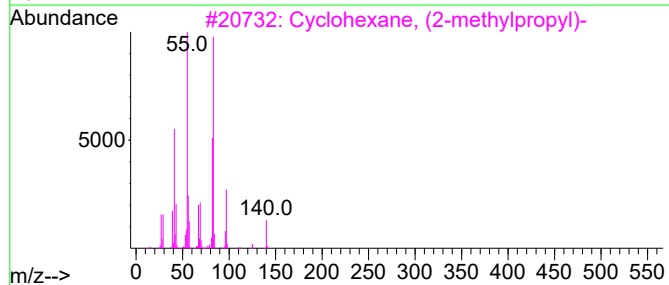
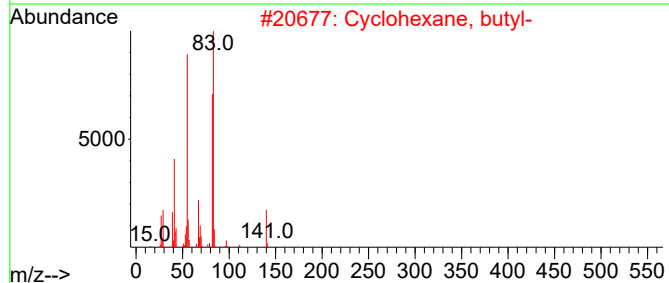
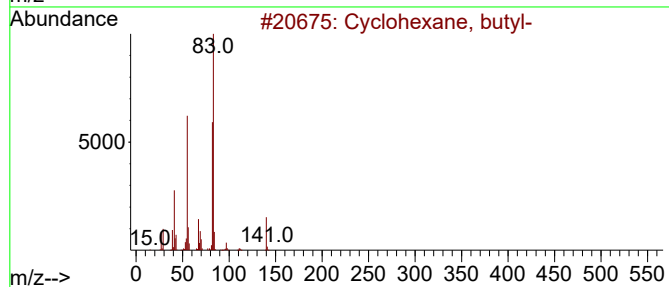
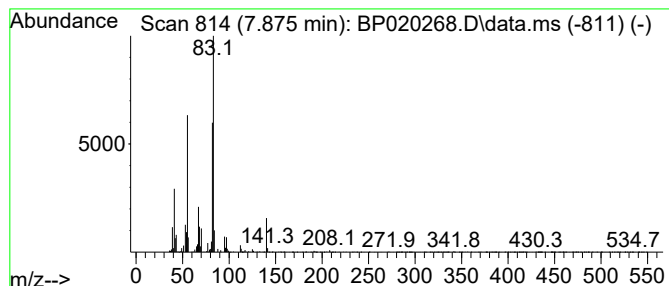
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Cyclic7.875 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.875	2.28 ng/ul	71226	1,4-Dichlorobenzene-d4	7.616	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, butyl-	140	C10H20	001678-93-9	90
2	Cyclohexane, butyl-	140	C10H20	001678-93-9	80
3	Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	72
4	Cyclohexane, butyl-	140	C10H20	001678-93-9	64
5	Cyclohexane, octyl-	196	C14H28	001795-15-9	64



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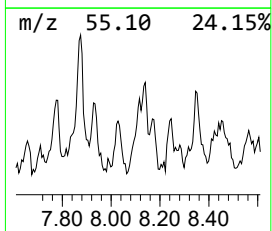
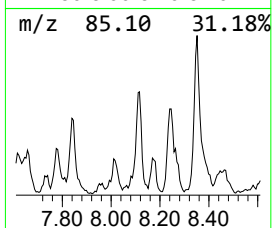
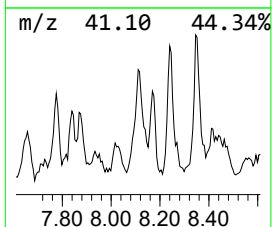
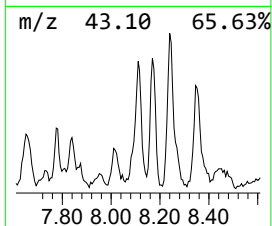
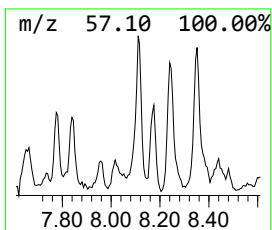
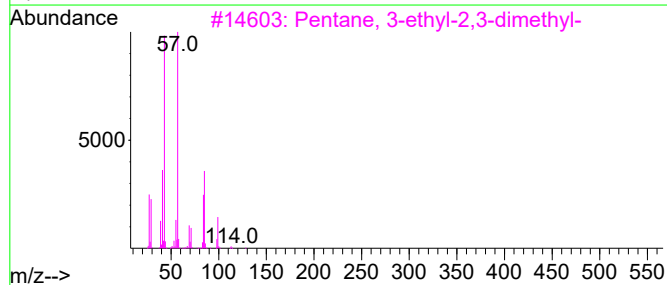
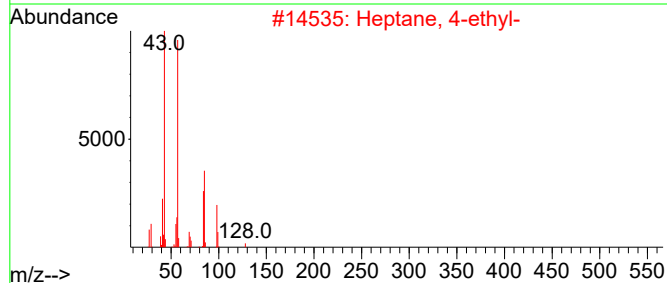
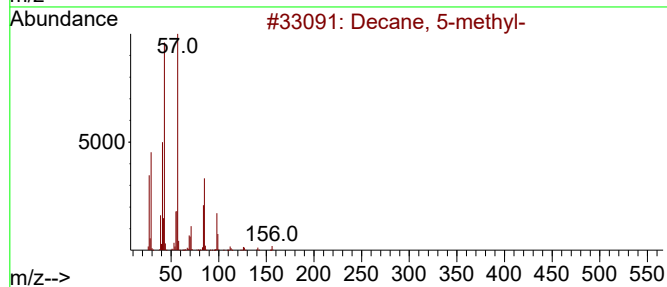
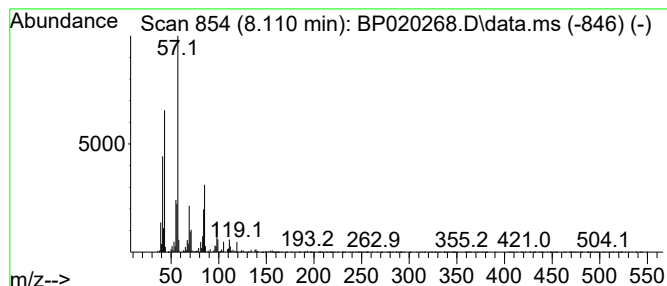
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.110	6.68 ng/ul	208958	1,4-Dichlorobenzene-d4			7.616
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane, 5-methyl-		156	C11H24	013151-35-4	59
2	Heptane, 4-ethyl-		128	C9H20	002216-32-2	53
3	Pentane, 3-ethyl-2,3-dimethyl-		128	C9H20	016747-33-4	50
4	Dodecane, 2,7,10-trimethyl-		212	C15H32	074645-98-0	50
5	Hexane, 2,4-dimethyl-		114	C8H18	000589-43-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
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Sample : P2369-16DL 30X
Misc :
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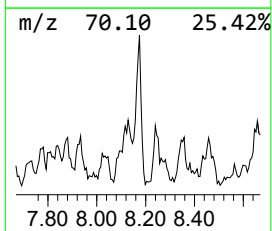
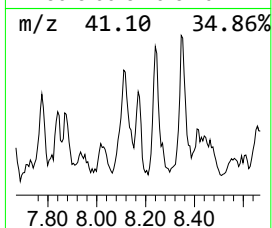
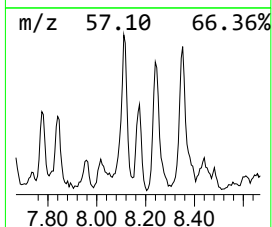
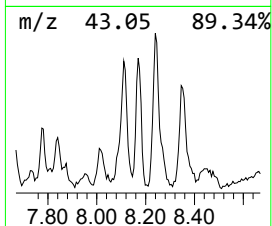
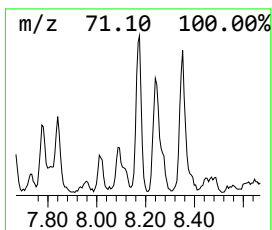
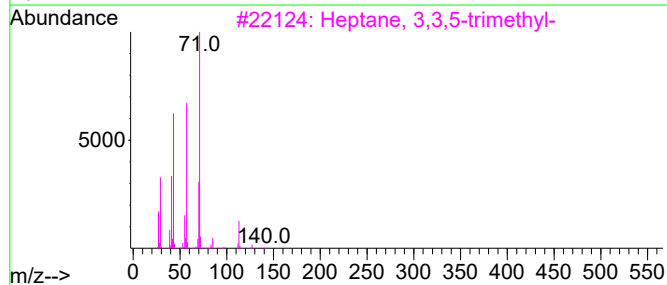
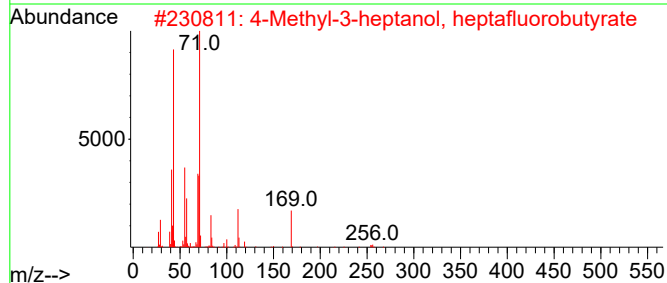
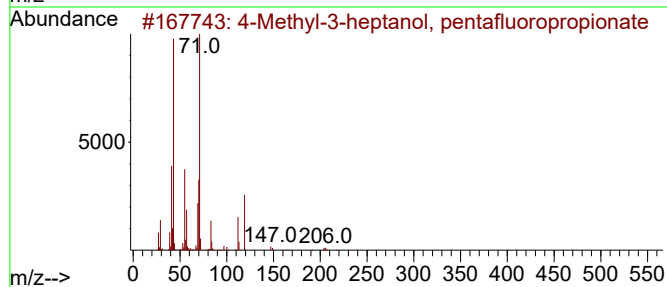
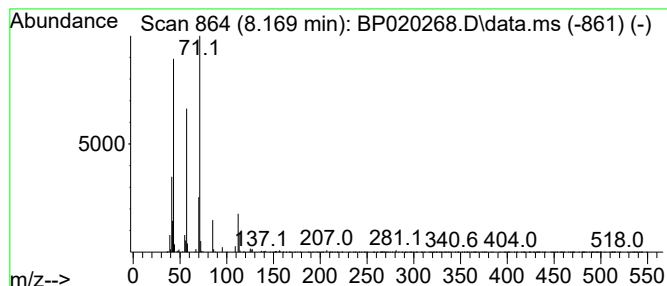
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 4-Methyl-3-heptanol, pentafl... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.169	3.84 ng/ul	120246	1,4-Dichlorobenzene-d4			7.616
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4-Methyl-3-heptanol, pentafluoro...		276	C11H17F5O2	1000365-51-7	78
2	4-Methyl-3-heptanol, heptafluoro...		326	C12H17F7O2	1000365-51-5	78
3	Heptane, 3,3,5-trimethyl-		142	C10H22	007154-80-5	78
4	Heptane, 2,5,5-trimethyl-		142	C10H22	001189-99-7	78
5	Hexane, 3,3,4,4-tetramethyl-		142	C10H22	005171-84-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
Operator : MA/JU
Sample : P2369-16DL 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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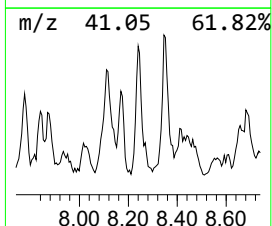
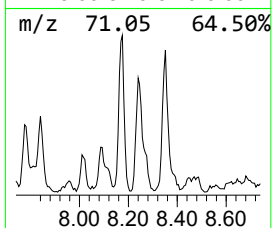
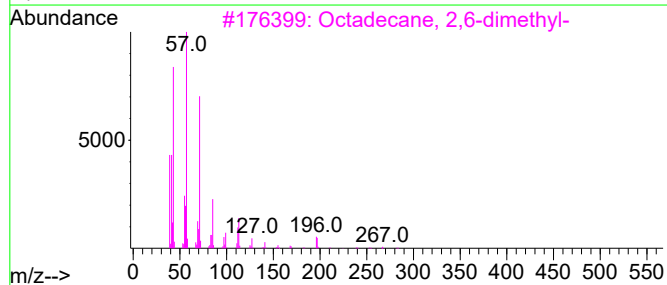
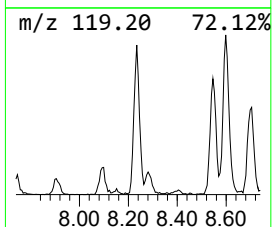
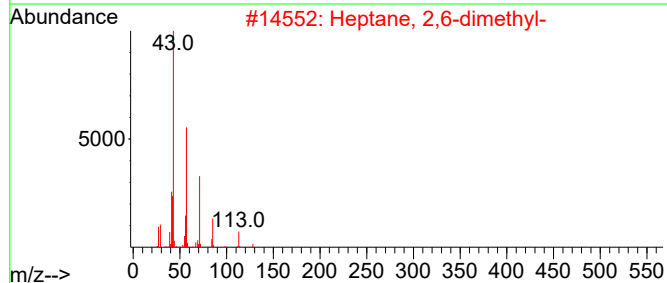
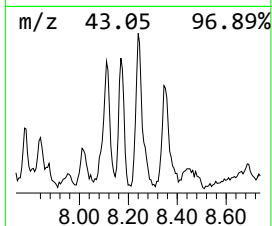
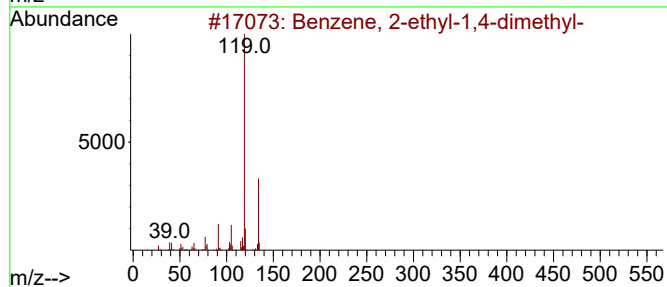
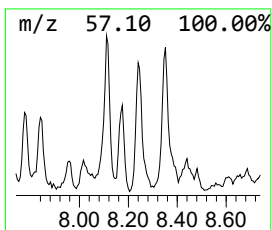
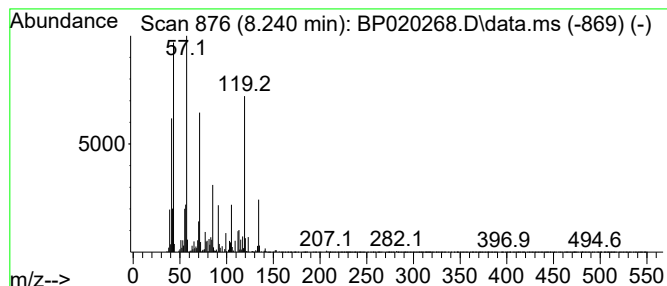
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.240	7.72 ng/ul	241679	1,4-Dichlorobenzene-d4	7.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	64
2			Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	60
3			Octadecane, 2,6-dimethyl-	282	C20H42	075163-97-2	55
4			Tridecane, 7-methyl-	198	C14H30	026730-14-3	50
5			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	44



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
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Sample : P2369-16DL 30X
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ALS Vial : 36 Sample Multiplier: 1

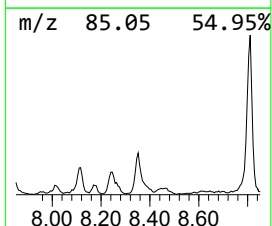
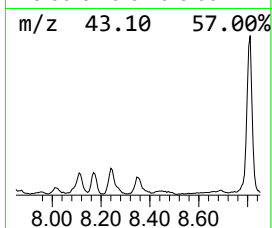
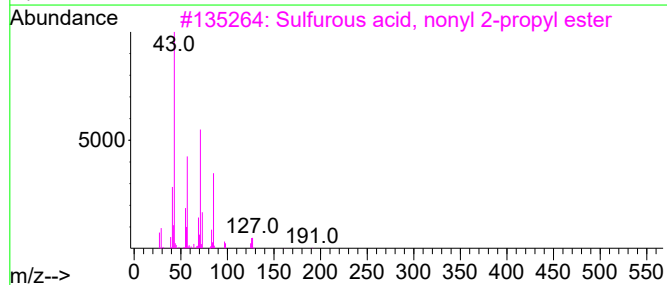
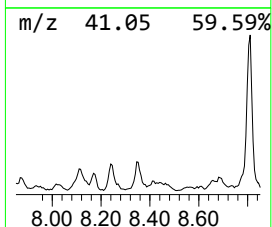
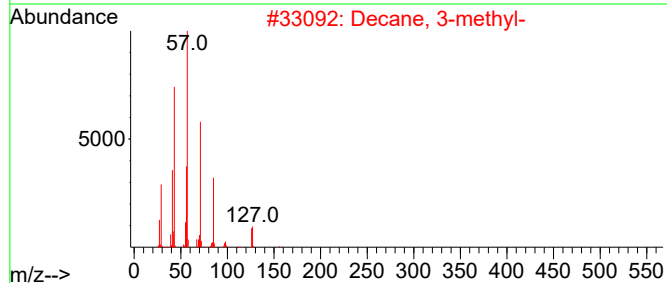
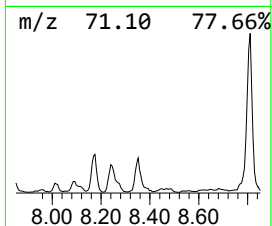
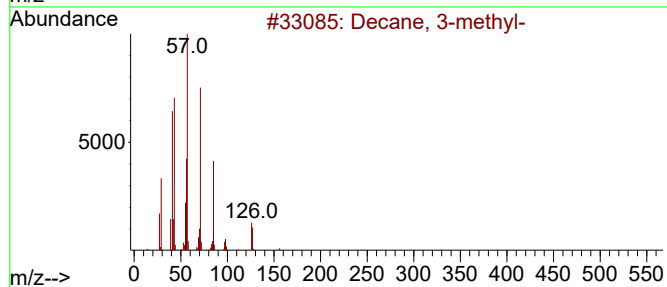
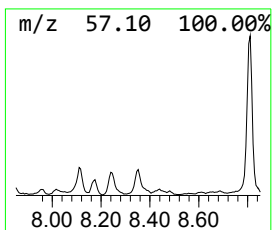
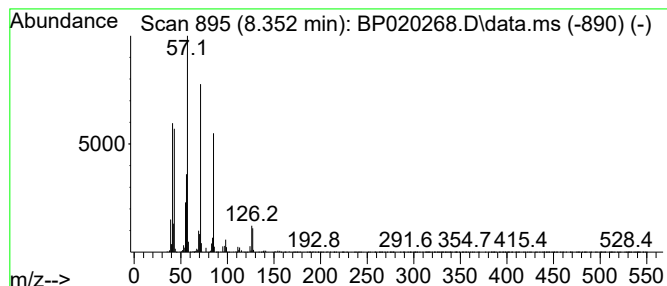
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.352	4.63 ng/ul	145012	1,4-Dichlorobenzene-d4			7.616
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane, 3-methyl-		156	C11H24	013151-34-3	91
2	Decane, 3-methyl-		156	C11H24	013151-34-3	90
3	Sulfurous acid, nonyl 2-propyl e...		250	C12H26O3S	1000309-12-0	72
4	Octane, 2-methyl-		128	C9H20	003221-61-2	59
5	Nonane, 1-iodo-		254	C9H19I	004282-42-2	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
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Sample : P2369-16DL 30X
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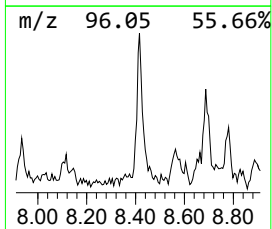
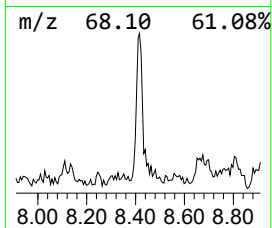
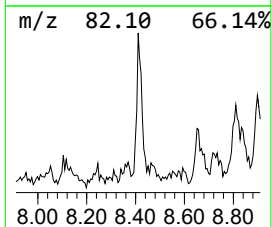
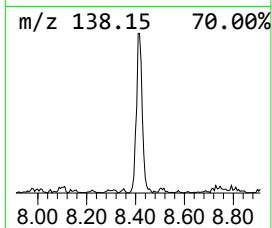
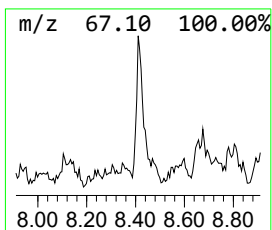
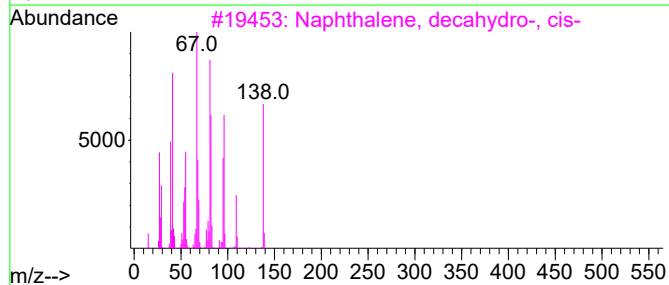
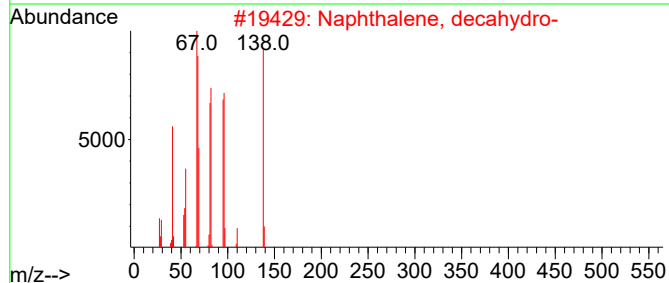
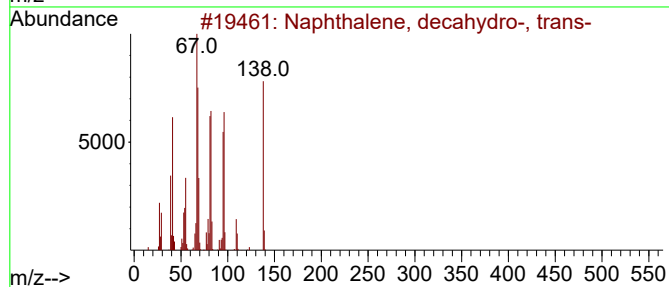
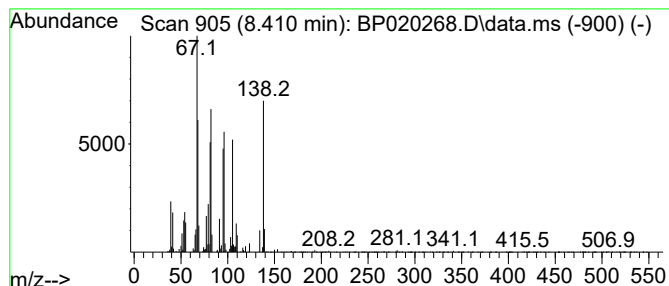
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Naphthalene, decahydro-, tr... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.410	9.44 ng/ul	295382	1,4-Dichlorobenzene-d4			7.616
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, decahydro-, trans-		138	C10H18	000493-02-7	94
2	Naphthalene, decahydro-		138	C10H18	000091-17-8	93
3	Naphthalene, decahydro-, cis-		138	C10H18	000493-01-6	83
4	Naphthalene, decahydro-, trans-		138	C10H18	000493-02-7	70
5	Bicyclo[2.2.1]heptane, 2-methyl-...		110	C8H14	000872-78-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
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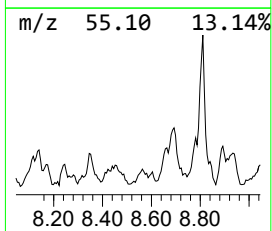
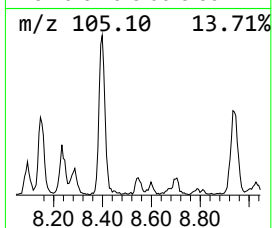
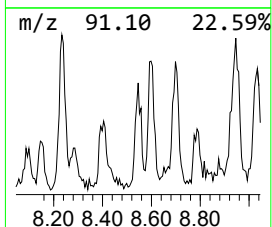
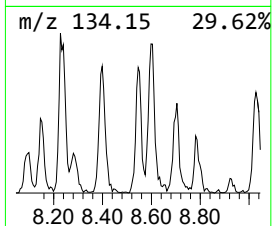
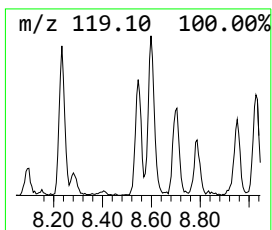
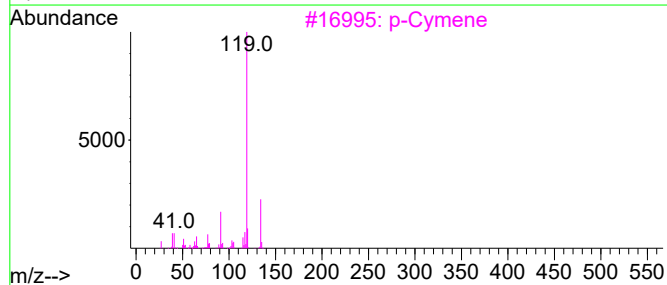
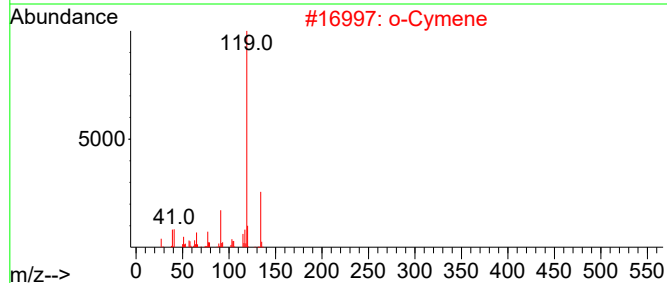
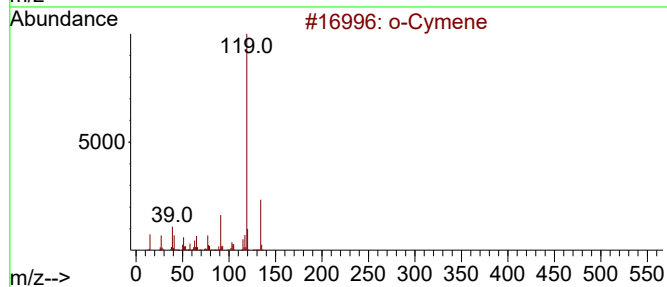
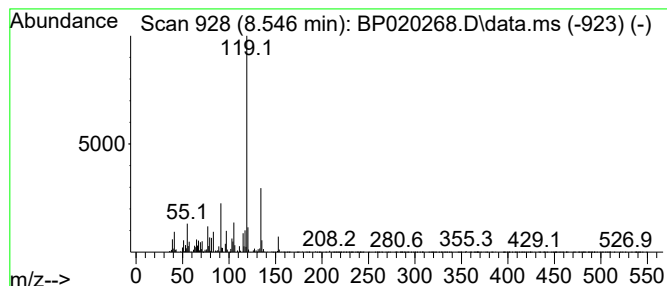
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 o-Cymene Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.546	2.57 ng/ul	80316	1,4-Dichlorobenzene-d4	7.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	95
2			o-Cymene	134	C10H14	000527-84-4	95
3			p-Cymene	134	C10H14	000099-87-6	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\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
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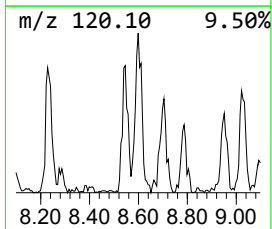
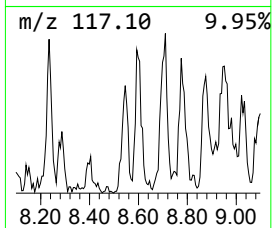
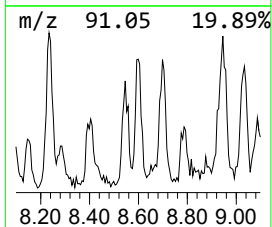
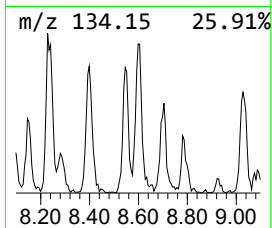
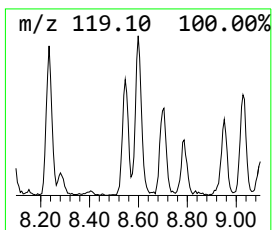
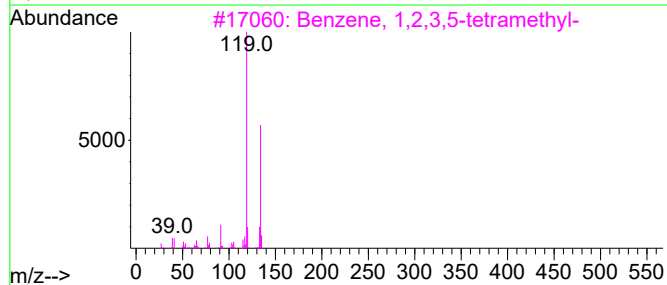
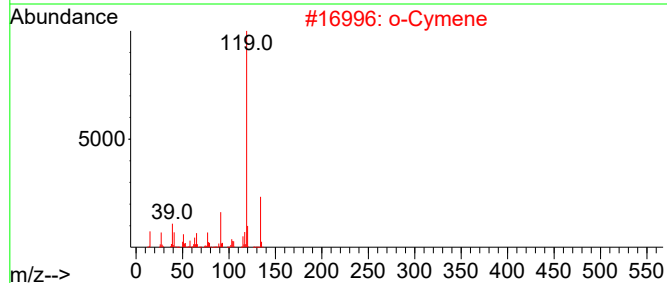
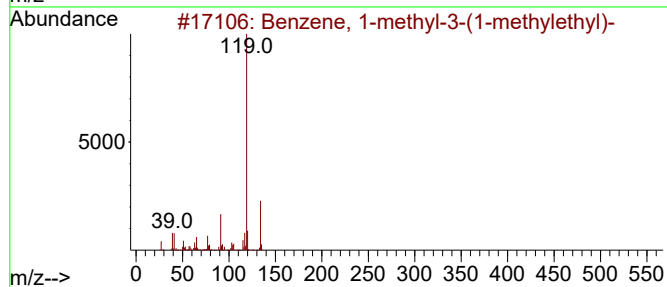
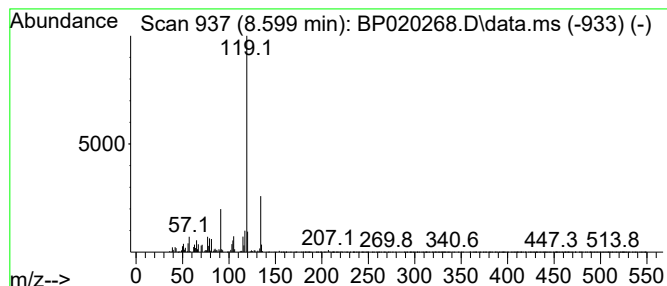
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Benzene, 1-methyl-3-(1-meth... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.599	2.20 ng/ul	68905	1,4-Dichlorobenzene-d4	7.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
2			o-Cymene	134	C10H14	000527-84-4	91
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
4			o-Cymene	134	C10H14	000527-84-4	91
5			o-Cymene	134	C10H14	000527-84-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
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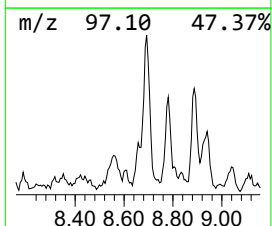
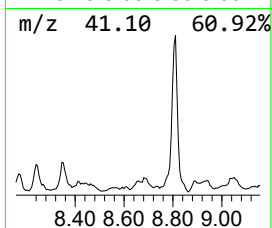
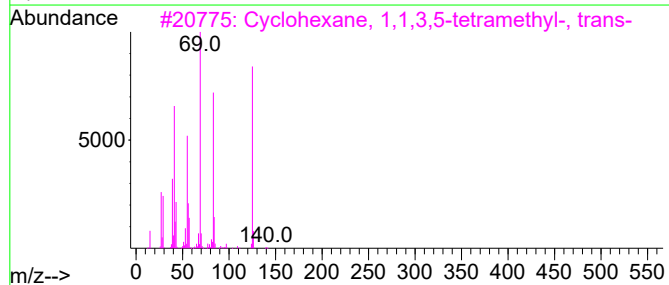
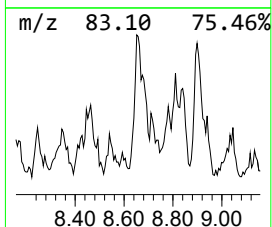
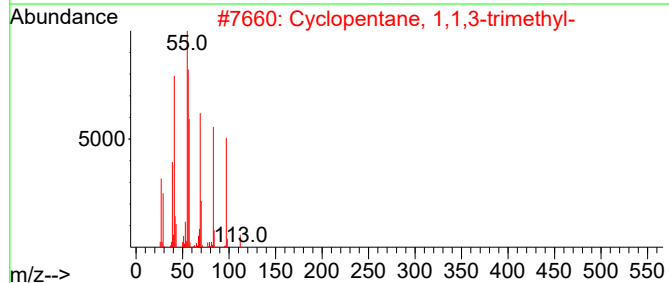
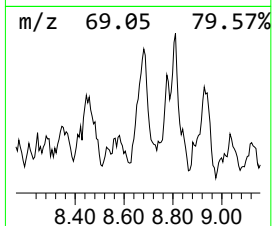
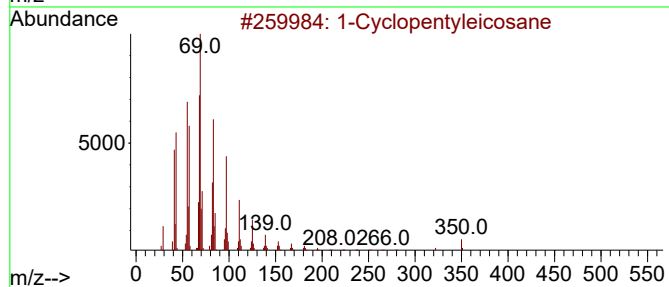
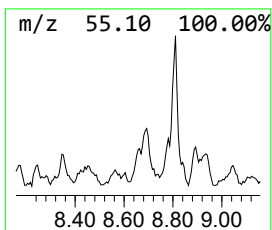
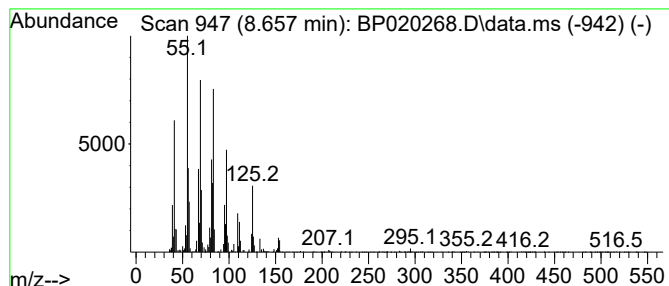
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Cyclic8.657 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.657	2.37 ng/ul	74080	1,4-Dichlorobenzene-d4	7.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Cyclopentyleicosane	350	C25H50	1000357-25-6	58
2			Cyclopentane, 1,1,3-trimethyl-	112	C8H16	004516-69-2	53
3			Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-31-8	43
4			Cyclohexane, 1,1,2-trimethyl-	126	C9H18	007094-26-0	38
5			Cyclopentane, 1-butyl-2-propyl-	168	C12H24	062199-50-2	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
Operator : MA/JU
Sample : P2369-16DL 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

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ClientSampleId :
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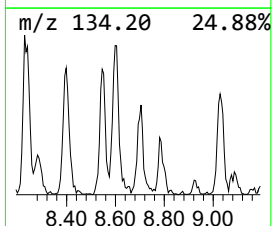
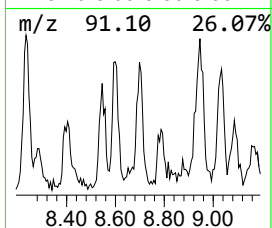
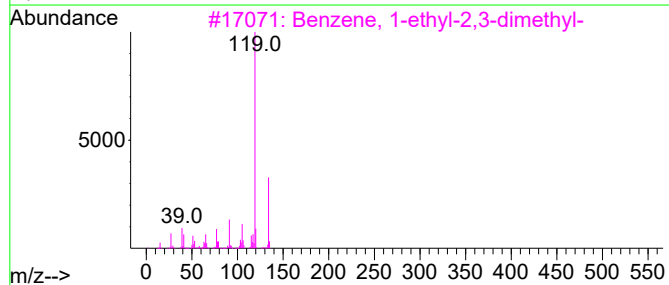
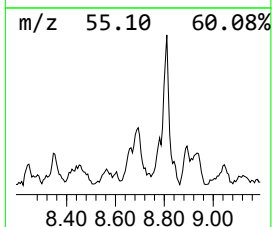
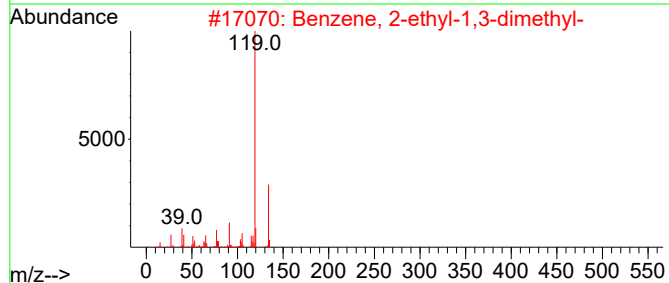
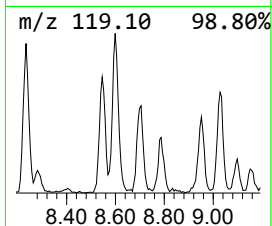
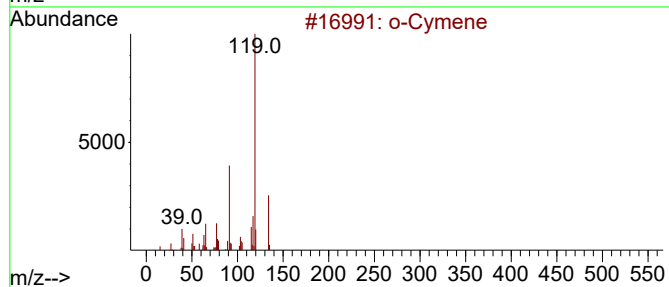
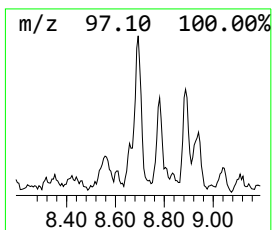
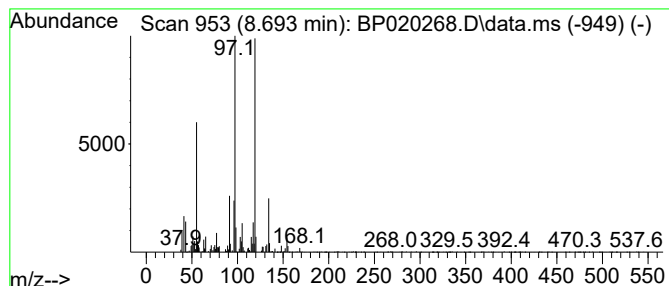
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.693	4.51 ng/ul	141276	1,4-Dichlorobenzene-d4	7.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	84
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	83
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	46
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	46
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
Operator : MA/JU
Sample : P2369-16DL 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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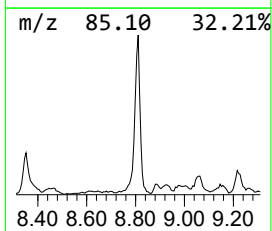
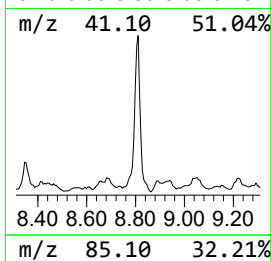
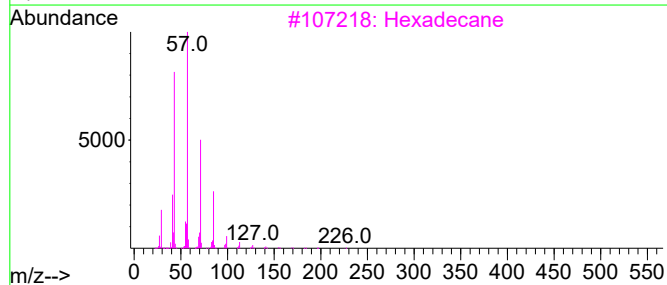
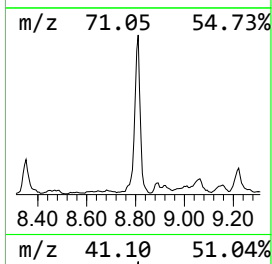
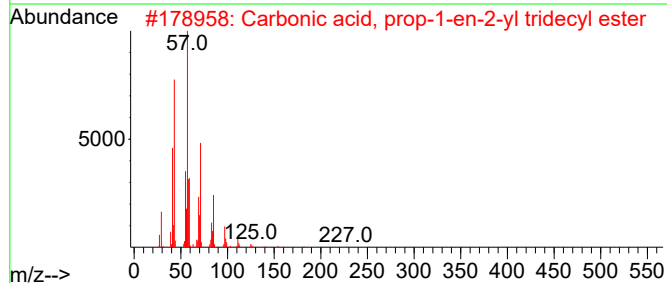
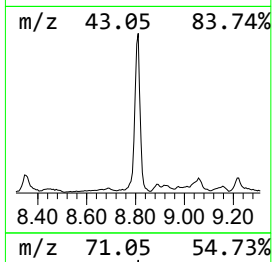
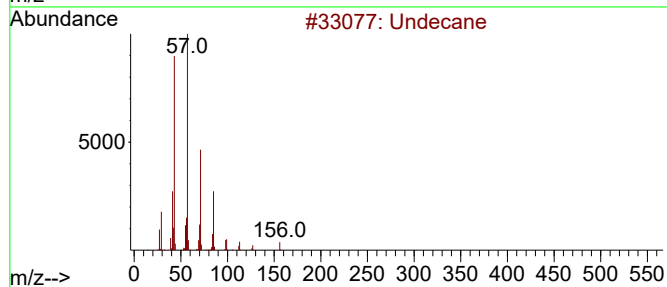
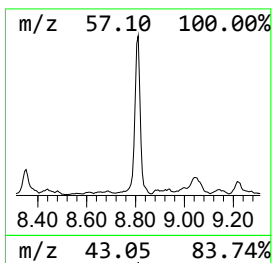
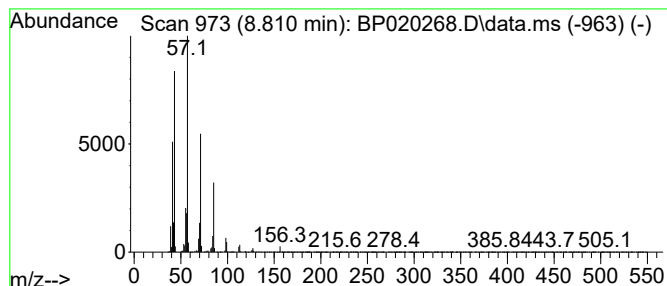
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.810	31.87 ng/ul	997659	1,4-Dichlorobenzene-d4	7.616

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 tr...	284	C17H32O3	1000382-90-8	90
3			Hexadecane	226	C16H34	000544-76-3	86
4			Decane, 6-ethyl-2-methyl-	184	C13H28	062108-21-8	83
5			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
Operator : MA/JU
Sample : P2369-16DL 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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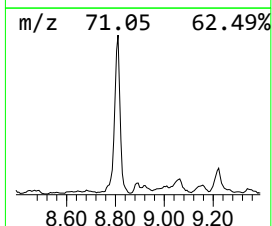
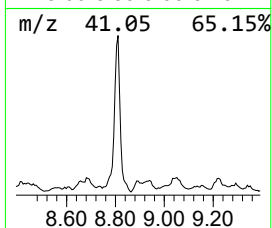
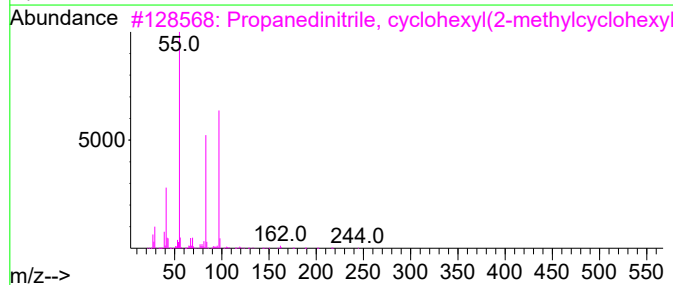
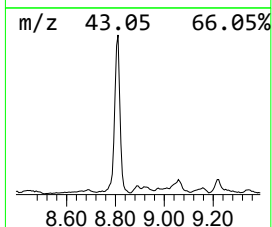
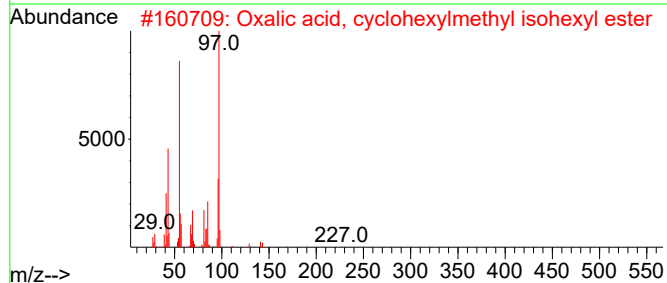
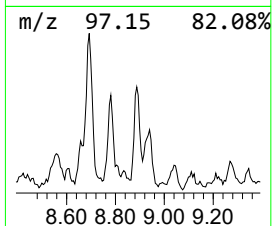
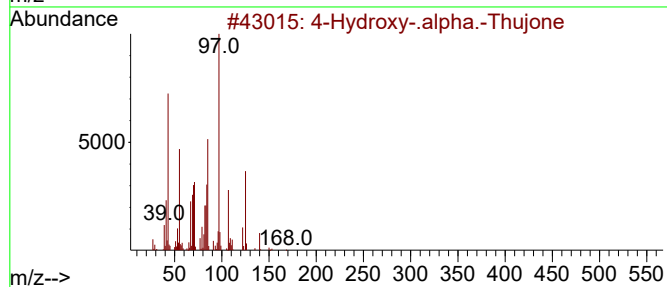
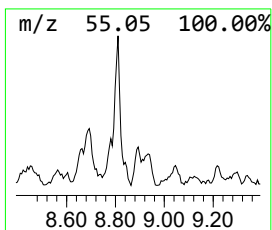
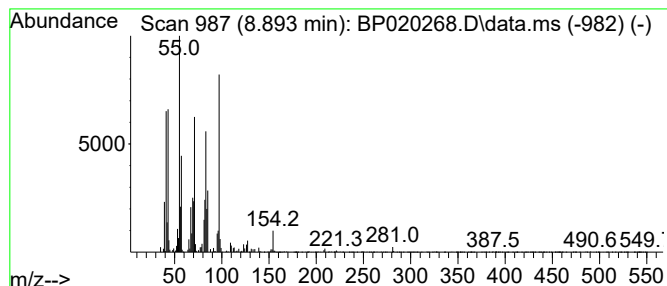
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 4-Hydroxy-.alpha.-Thujone Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.893	2.88 ng/ul	90184	1,4-Dichlorobenzene-d4	7.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Hydroxy-.alpha.-Thujone	168	C10H16O2	158612-68-1	52
2			Oxalic acid, cyclohexylmethyl is...	270	C15H26O4	1000309-68-3	43
3			Propanedinitrile, cyclohexyl(2-m...	244	C16H24N2	074764-55-9	35
4			Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	27
5			Cyclopentane, 1-hexyl-3-methyl-	168	C12H24	061142-68-5	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
Operator : MA/JU
Sample : P2369-16DL 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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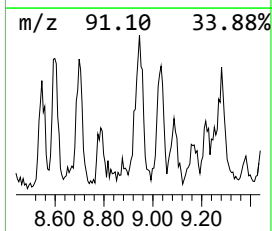
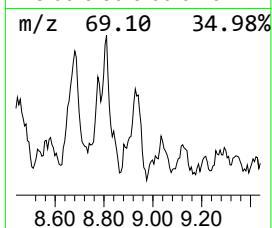
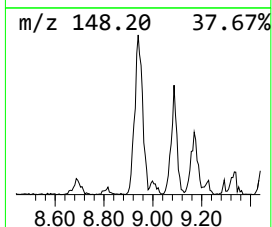
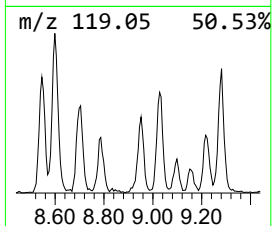
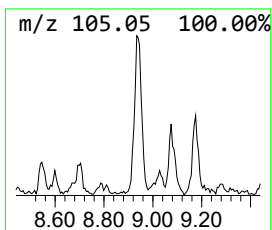
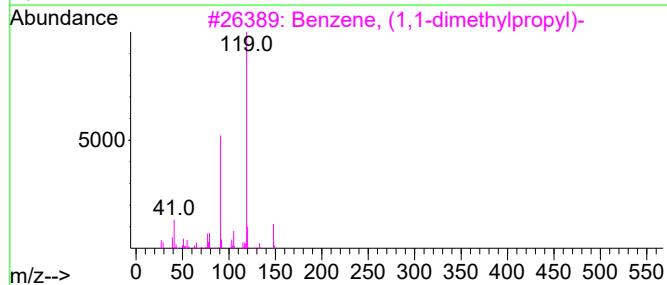
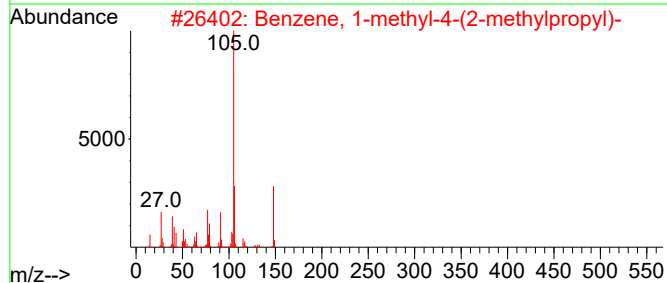
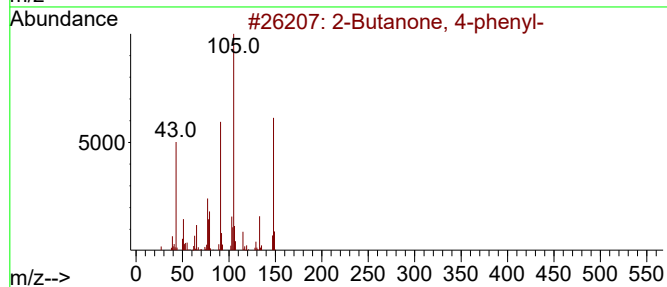
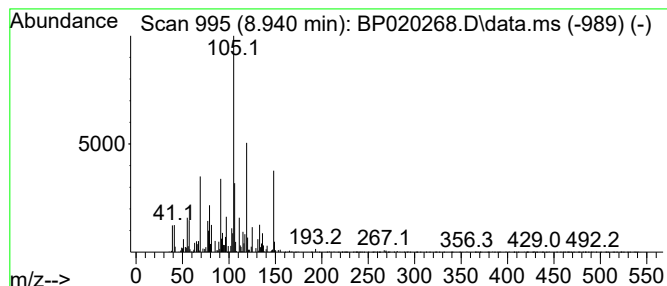
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.940	5.52 ng/ul	172847	1,4-Dichlorobenzene-d4	7.616

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Butanone, 4-phenyl-	148	C10H12O	002550-26-7	42
2		Benzene, 1-methyl-4-(2-methylpropyl)-	148	C11H16	005161-04-6	38
3		Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	35
4		Benzene, (1,2-dimethylpropyl)-	148	C11H16	004481-30-5	30
5		2-Butanone, 4-phenyl-	148	C10H12O	002550-26-7	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
Operator : MA/JU
Sample : P2369-16DL 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
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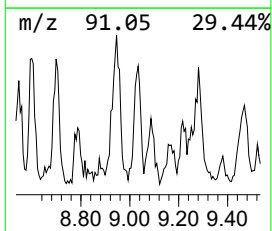
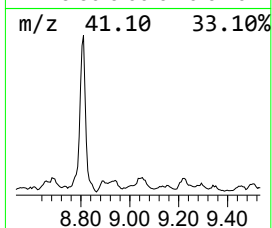
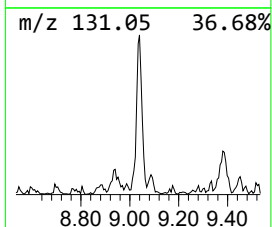
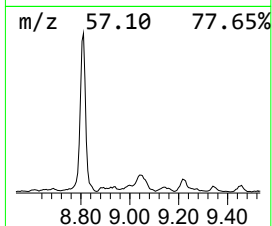
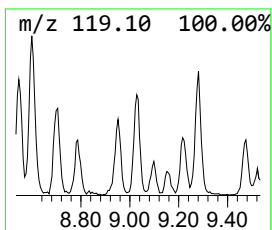
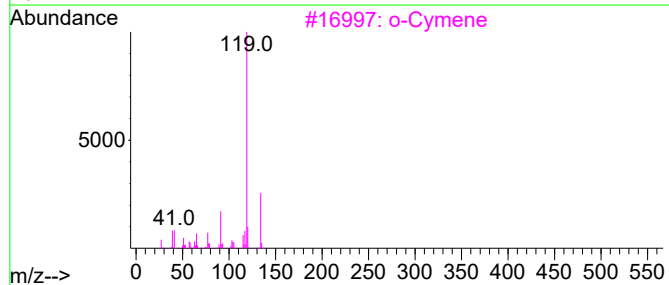
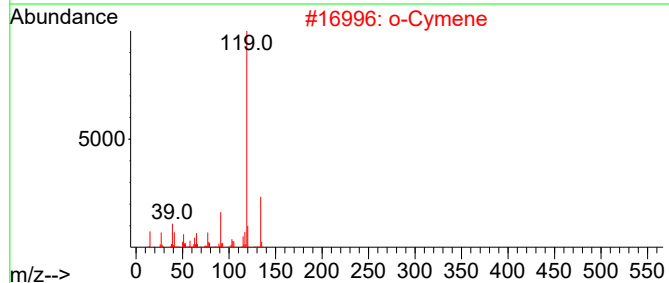
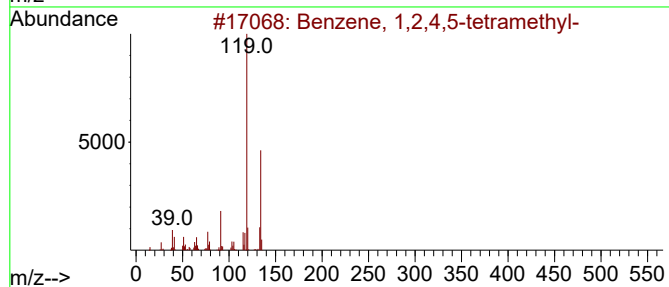
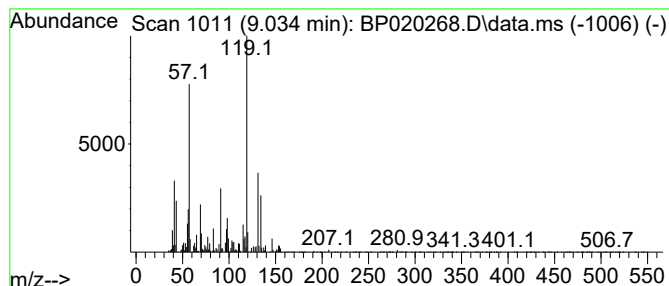
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.034	2.59 ng/ul	116562	Naphthalene-d8	10.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	55
2			o-Cymene	134	C10H14	000527-84-4	55
3			o-Cymene	134	C10H14	000527-84-4	55
4			p-Cymene	134	C10H14	000099-87-6	49
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
Operator : MA/JU
Sample : P2369-16DL 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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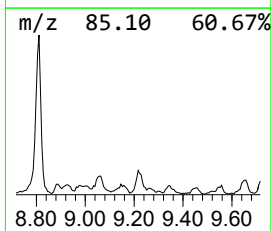
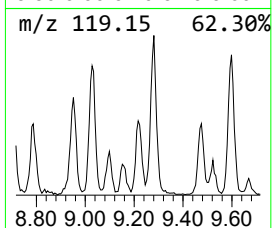
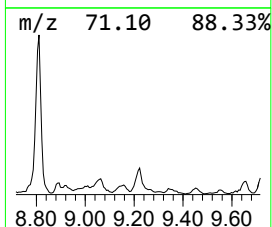
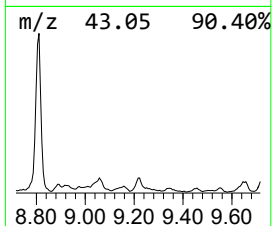
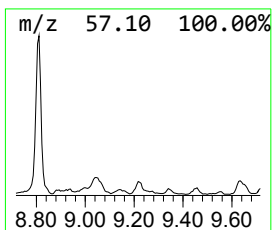
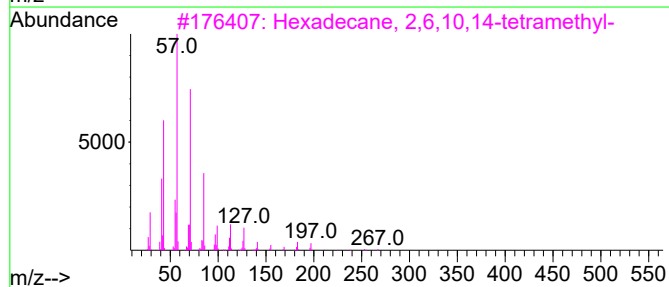
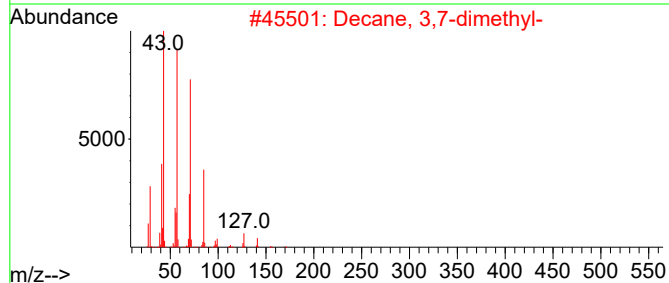
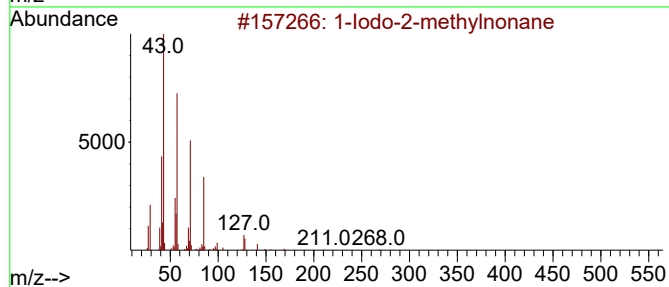
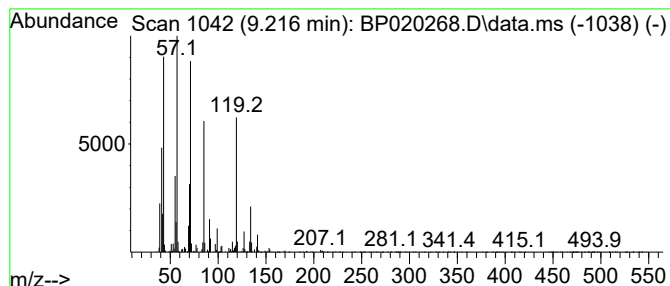
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 1-Iodo-2-methylnonane Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.216	2.25 ng/ul	101215	Naphthalene-d8	10.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	64
2			Decane, 3,7-dimethyl-	170	C12H26	017312-54-8	62
3			Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	53
4			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	52
5			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
Operator : MA/JU
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Misc :
ALS Vial : 36 Sample Multiplier: 1

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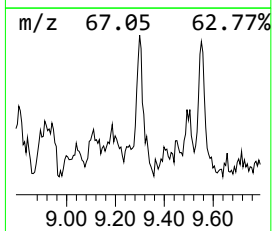
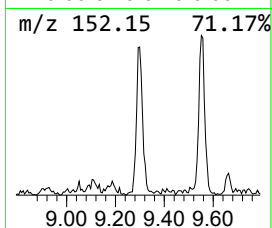
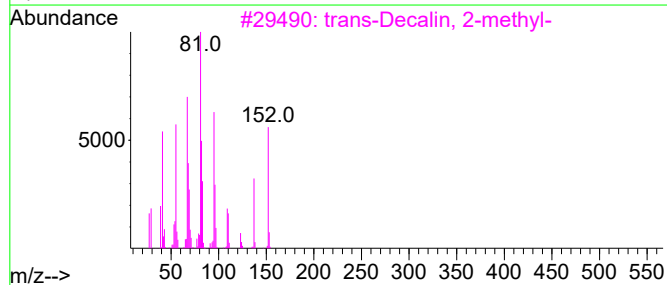
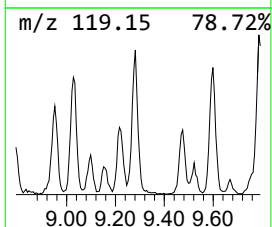
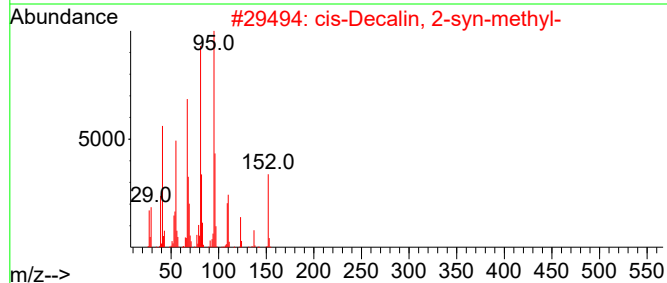
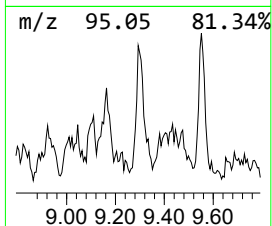
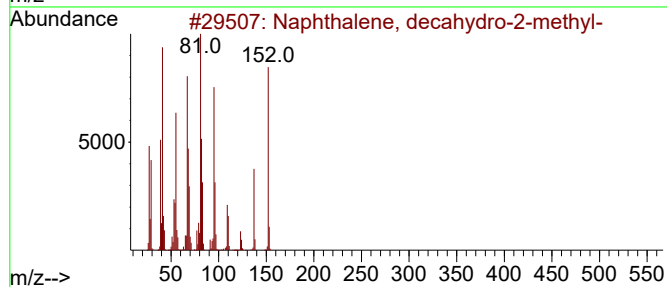
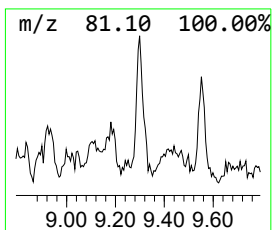
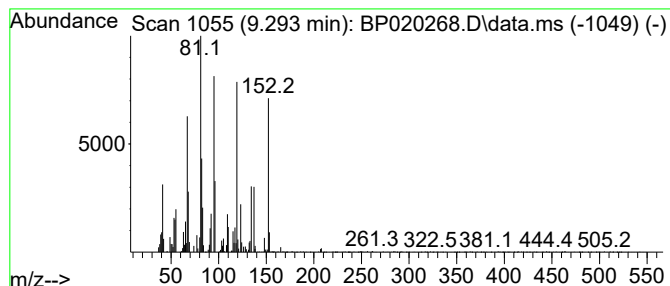
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Naphthalene, decahydro-2-me... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.293	2.26 ng/ul	101688	Naphthalene-d8	10.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	94
2			cis-Decalin, 2-syn-methyl-	152	C11H20	1000155-85-6	64
3			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	58
4			Pulegone	152	C10H16O	000089-82-7	50
5			Pulegone	152	C10H16O	000089-82-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
Operator : MA/JU
Sample : P2369-16DL 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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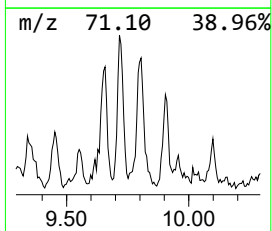
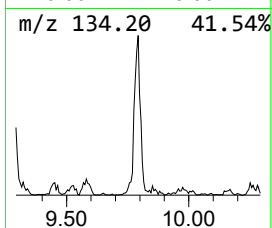
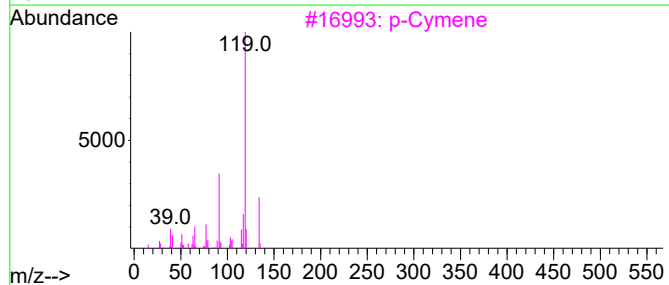
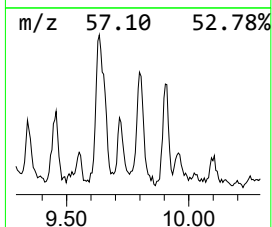
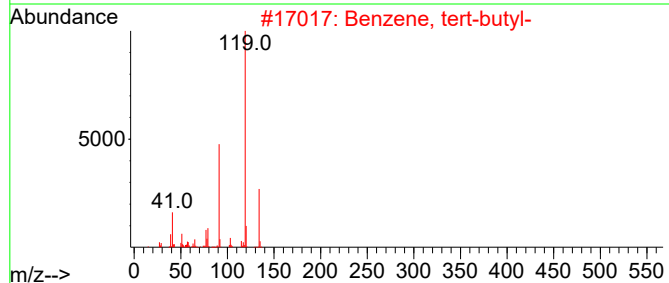
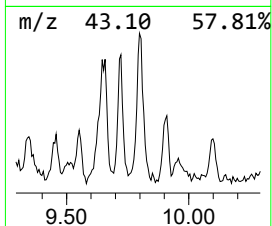
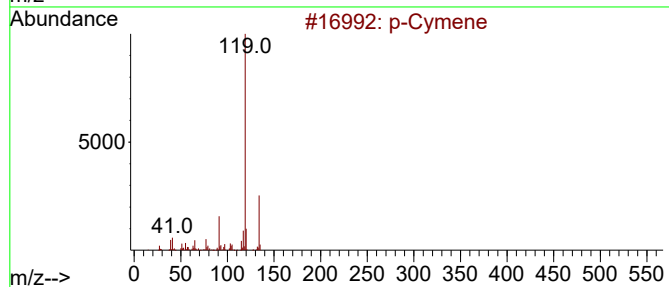
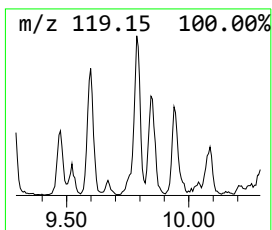
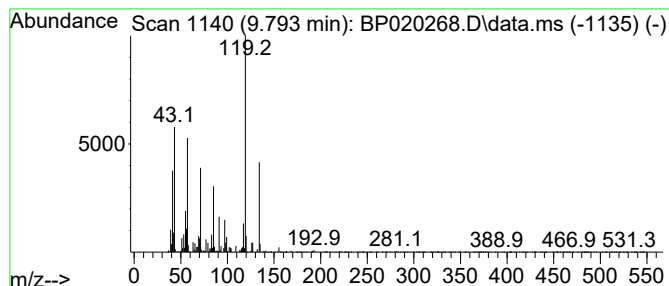
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 p-Cymene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.793	3.35 ng/ul	150810	Naphthalene-d8	10.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	64
2			Benzene, tert-butyl-	134	C10H14	000098-06-6	60
3			p-Cymene	134	C10H14	000099-87-6	53
4			o-Cymene	134	C10H14	000527-84-4	49
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
Operator : MA/JU
Sample : P2369-16DL 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43P0DL

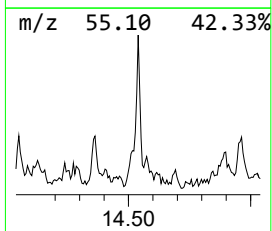
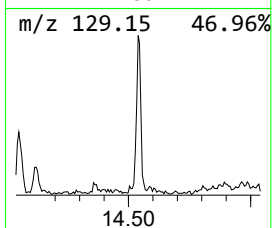
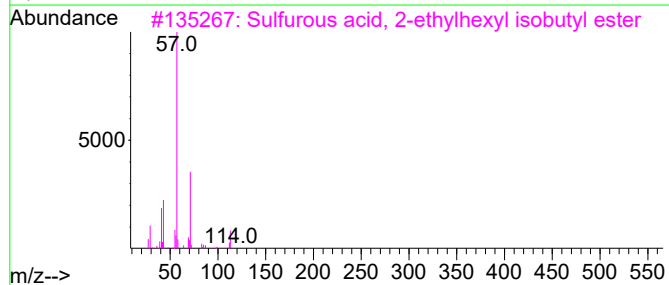
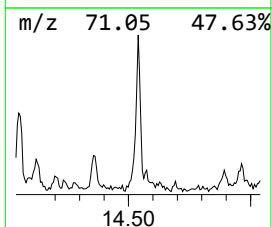
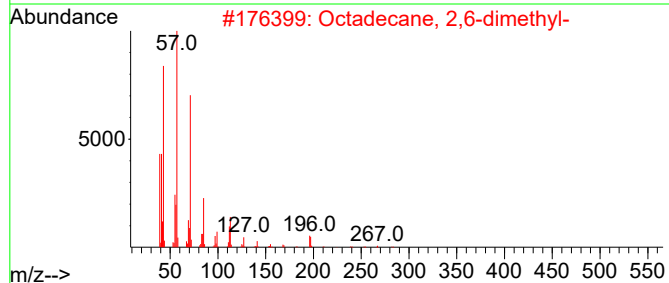
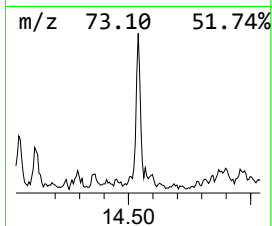
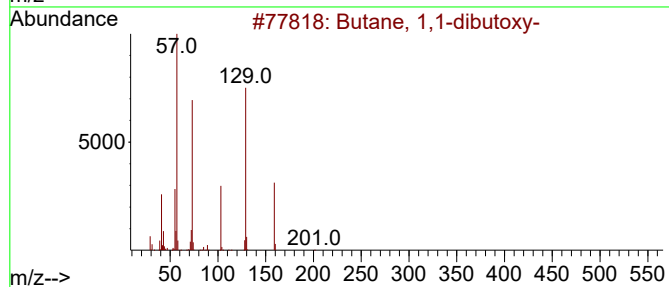
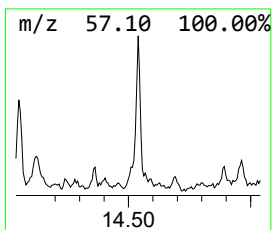
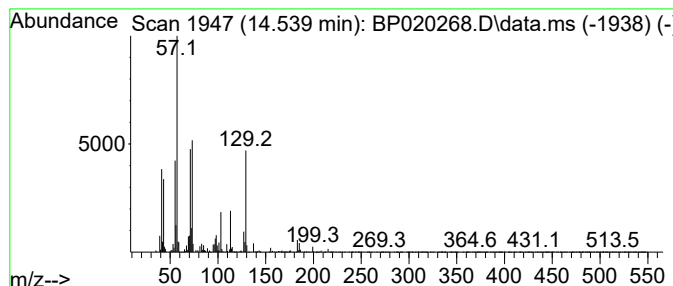
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 unknown-02 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.540	2.54 ng/ul	153943	Acenaphthene-d10	14.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 1,1-dibutoxy-	202	C12H26O2	005921-80-2	43		
2	Octadecane, 2,6-dimethyl-	282	C20H42	075163-97-2	35		
3	Sulfurous acid, 2-ethylhexyl iso...	250	C12H26O3S	1000309-18-7	32		
4	Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	32		
5	Sulfurous acid, di(2-ethylhexyl)...	306	C16H34O3S	1000309-19-1	32		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020268.D
Acq On : 09 May 2024 13:42
Operator : MA/JU
Sample : P2369-16DL 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43P0DL

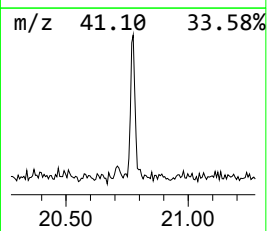
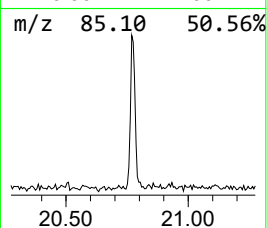
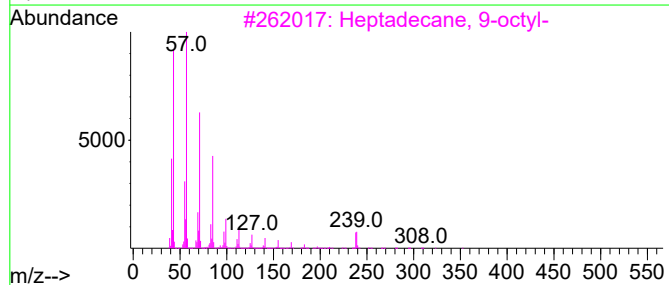
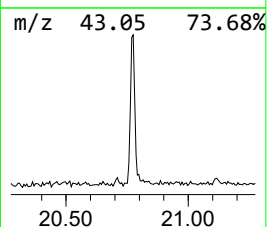
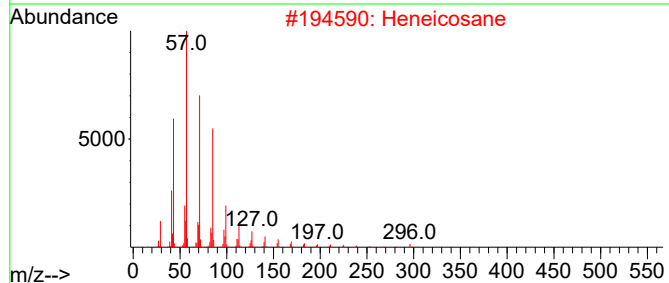
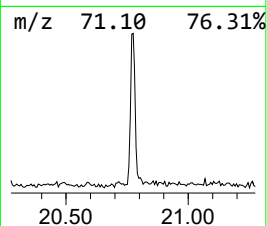
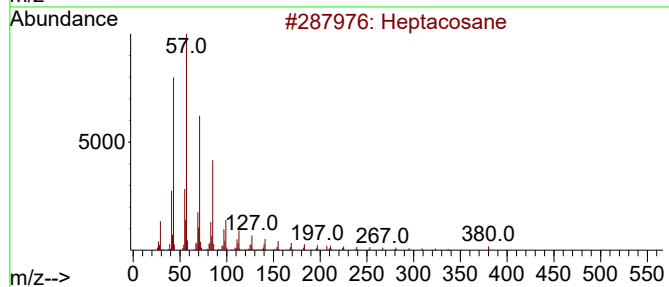
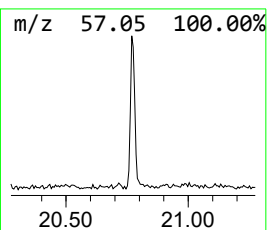
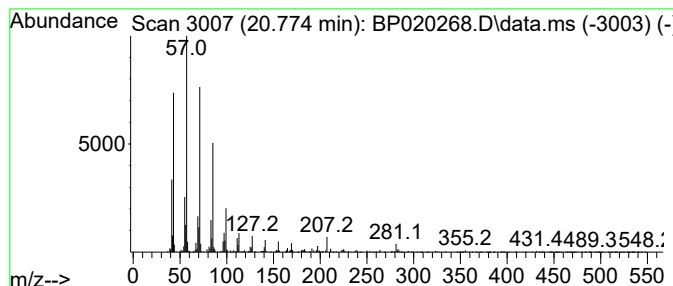
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.774	2.09 ng/ul	135124	Chrysene-d12	21.633

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptacosane	380	C27H56	000593-49-7	87
2		Heneicosane	296	C21H44	000629-94-7	87
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	86
4		Octacosane	394	C28H58	000630-02-4	81
5		Tridecane, 6-propyl-	226	C16H34	055045-10-8	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
 Data File : BP020268.D
 Acq On : 09 May 2024 13:42
 Operator : MA/JU
 Sample : P2369-16DL 30X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43P0DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	6.275	2.0	ng/ul	63809	1	7.616	626050	20.0
(DEL) Alkane: S...	6.763	2.6	ng/ul	81741	1	7.616	626050	20.0
(DEL) Alkane: S...	7.222	7.8	ng/ul	243050	1	7.616	626050	20.0
Hexadecane, 1-c...	7.510	4.4	ng/ul	136406	1	7.616	626050	20.0
(DEL) Alkane: S...	7.563	7.0	ng/ul	219039	1	7.616	626050	20.0
Benzene, 1,2,3-...	7.722	3.1	ng/ul	95925	1	7.616	626050	20.0
(DEL) Alkane: S...	7.775	2.9	ng/ul	91331	1	7.616	626050	20.0
(DEL) Alkane: C...	7.875	2.3	ng/ul	71226	1	7.616	626050	20.0
(DEL) Alkane: S...	8.110	6.7	ng/ul	208958	1	7.616	626050	20.0
4-Methyl-3-hept...	8.169	3.8	ng/ul	120246	1	7.616	626050	20.0
Benzene, 2-ethy...	8.240	7.7	ng/ul	241679	1	7.616	626050	20.0
(DEL) Alkane: S...	8.352	4.6	ng/ul	145012	1	7.616	626050	20.0
Naphthalene, de...	8.410	9.4	ng/ul	295382	1	7.616	626050	20.0
o-Cymene	8.546	2.6	ng/ul	80316	1	7.616	626050	20.0
Benzene, 1-meth...	8.599	2.2	ng/ul	68905	1	7.616	626050	20.0
(DEL) Alkane: C...	8.657	2.4	ng/ul	74080	1	7.616	626050	20.0
Benzene, 2-ethy...	8.693	4.5	ng/ul	141276	1	7.616	626050	20.0
(DEL) Alkane: S...	8.810	31.9	ng/ul	997659	1	7.616	626050	20.0
4-Hydroxy-.alph...	8.893	2.9	ng/ul	90184	1	7.616	626050	20.0
unknown-01	8.940	5.5	ng/ul	172847	1	7.616	626050	20.0
Benzene, 1,2,4,...	9.034	2.6	ng/ul	116562	2	10.387	899073	20.0
1-Iodo-2-methyl...	9.216	2.3	ng/ul	101215	2	10.387	899073	20.0
Naphthalene, de...	9.293	2.3	ng/ul	101688	2	10.387	899073	20.0
p-Cymene	9.793	3.4	ng/ul	150810	2	10.387	899073	20.0
unknown-02	14.540	2.5	ng/ul	153943	3	14.292	1211510	20.0
(DEL) Alkane: S...	20.774	2.1	ng/ul	135124	5	21.633	1290160	20.0