

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
 Data File : BP013279.D
 Acq On : 23 Dec 2022 14:52
 Operator : CG/JU
 Sample : N6018-09
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBYH9

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-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP013279.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.334	21	25	36	rVB	128108	178388	1.97%	0.130%
2	3.764	94	98	113	rBV	1609358	2470334	27.32%	1.799%
3	4.087	149	153	159	rBV	45199	61263	0.68%	0.045%
4	4.464	213	217	222	rVB7	14024	21883	0.24%	0.016%
5	5.034	310	314	320	rVB2	58477	87889	0.97%	0.064%
6	5.263	347	353	359	rBV	53404	94276	1.04%	0.069%
7	6.458	550	556	563	rBV2	23338	38348	0.42%	0.028%
8	6.528	565	568	575	rVV6	15164	29276	0.32%	0.021%
9	6.616	577	583	590	rVB	263447	415100	4.59%	0.302%
10	6.922	629	635	643	rBV	122011	193389	2.14%	0.141%
11	7.105	659	666	676	rVV	2135372	3354380	37.10%	2.443%
12	7.269	685	694	705	rVB	1759612	2640464	29.20%	1.923%
13	7.375	706	712	718	rVB	325987	493583	5.46%	0.359%
14	7.475	721	729	739	rBV	2168680	3447101	38.12%	2.511%
15	7.946	798	809	816	rBV	2356491	3732839	41.28%	2.719%
16	8.005	816	819	825	rVB5	14648	27628	0.31%	0.020%
17	8.075	825	831	835	rBV3	14005	29442	0.33%	0.021%
18	8.157	840	845	850	rBV5	24134	42203	0.47%	0.031%
19	8.657	923	930	938	rVV	2368806	3891579	43.04%	2.834%
20	9.110	1000	1007	1017	rVV	2019490	3339749	36.93%	2.432%
21	9.269	1029	1034	1041	rVB4	26564	42464	0.47%	0.031%
22	9.510	1070	1075	1081	rBV2	56982	101156	1.12%	0.074%
23	9.699	1103	1107	1115	rVB	16159	37350	0.41%	0.027%
24	9.840	1121	1131	1144	rBV	1699568	2895101	32.02%	2.109%
25	10.063	1161	1169	1175	rVB10	29352	59560	0.66%	0.043%
26	10.263	1196	1203	1209	rBV4	15736	28206	0.31%	0.021%
27	10.375	1215	1222	1234	rBV	2672765	4508921	49.86%	3.284%
28	10.534	1244	1249	1256	rVB7	15237	30264	0.33%	0.022%
29	10.757	1279	1287	1295	rBV	3283558	5471640	60.51%	3.985%
30	10.898	1302	1311	1330	rBV	1987392	3503900	38.75%	2.552%
31	12.163	1523	1526	1535	rVB3	21100	33141	0.37%	0.024%
32	12.240	1535	1539	1547	rBV2	16275	28975	0.32%	0.021%
33	12.340	1547	1556	1561	rBV	47803	101094	1.12%	0.074%
34	12.787	1627	1632	1636	rVV7	16474	31731	0.35%	0.023%
35	12.951	1655	1660	1661	rVV5	15561	23514	0.26%	0.017%
36	13.016	1667	1671	1676	rVB6	18373	27097	0.30%	0.020%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
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37	13.322	1718	1723	1725	rVV6	15366	25085	0.28%	0.018%
38	13.345	1725	1727	1729	rVV3	18203	21841	0.24%	0.016%
39	13.375	1729	1732	1733	rVV3	27684	33018	0.37%	0.024%
40	13.398	1733	1736	1740	rVB	50027	71875	0.79%	0.052%
41	13.457	1740	1746	1752	rBV	513135	717798	7.94%	0.523%
42	13.740	1791	1794	1802	rVB9	18348	25612	0.28%	0.019%
43	13.857	1806	1814	1818	rBV7	50783	90198	1.00%	0.066%
44	13.904	1818	1822	1829	rVB2	142397	209186	2.31%	0.152%
45	13.992	1829	1837	1846	rBV	4376231	6105818	67.53%	4.447%
46	14.057	1846	1848	1852	rVB4	23482	25898	0.29%	0.019%
47	14.139	1853	1862	1866	rBV	301468	455767	5.04%	0.332%
48	14.175	1866	1868	1874	rVB2	48024	77854	0.86%	0.057%
49	14.281	1876	1886	1894	rBV	5093629	7453941	82.43%	5.429%
50	14.422	1905	1910	1915	rBV2	77877	112257	1.24%	0.082%
51	14.475	1915	1919	1924	rVB2	53036	74251	0.82%	0.054%
52	14.539	1926	1930	1933	rBV	114097	163166	1.80%	0.119%
53	14.587	1933	1938	1944	rVV	4703072	6878281	76.07%	5.010%
54	14.639	1944	1947	1952	rVV2	309194	591510	6.54%	0.431%
55	14.687	1952	1955	1966	rVV5	129624	264847	2.93%	0.193%
56	14.792	1966	1973	1978	rVV	1242794	1997149	22.09%	1.455%
57	14.828	1978	1979	1984	rVV	229453	322080	3.56%	0.235%
58	14.892	1985	1990	1994	rVV	600802	1028807	11.38%	0.749%
59	14.951	1994	2000	2025	rVB3	810403	2628462	29.07%	1.914%
60	15.345	2063	2067	2075	rVB8	49851	93363	1.03%	0.068%
61	15.481	2085	2090	2099	rBV4	178467	419123	4.64%	0.305%
62	15.581	2101	2107	2122	rVV	6104634	8821448	97.56%	6.425%
63	15.704	2122	2128	2134	rVV	1227410	1633784	18.07%	1.190%
64	15.816	2143	2147	2153	rVB5	72782	116935	1.29%	0.085%
65	15.875	2154	2157	2161	rVB4	36678	45434	0.50%	0.033%
66	15.951	2164	2170	2177	rVB	772082	1052241	11.64%	0.766%
67	16.034	2180	2184	2185	rBV3	30807	36600	0.40%	0.027%
68	16.063	2185	2189	2198	rVB3	92329	181290	2.00%	0.132%
69	16.151	2198	2204	2209	rBV7	49418	85888	0.95%	0.063%
70	16.210	2211	2214	2217	rBV5	19789	29921	0.33%	0.022%
71	16.292	2225	2228	2231	rBV5	15863	26761	0.30%	0.019%
72	16.433	2248	2252	2258	rBV8	43005	84104	0.93%	0.061%
73	16.516	2261	2266	2270	rVV2	151317	205293	2.27%	0.150%
74	16.733	2299	2303	2307	rBV3	64982	99715	1.10%	0.073%
75	16.845	2319	2322	2327	rVB6	28018	32976	0.36%	0.024%
76	17.092	2361	2364	2367	rBV5	16794	22277	0.25%	0.016%

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77	17.128	2367	2370	2374	rVB5	26141	36715	0.41%	0.027%
78	17.228	2382	2387	2391	rBV2	84519	131159	1.45%	0.096%
79	17.298	2395	2399	2401	rBV3	51857	63350	0.70%	0.046%
80	17.345	2401	2407	2418	rVV	4647274	6983013	77.23%	5.086%
81	17.445	2418	2424	2433	rVV	5471545	7742814	85.63%	5.639%
82	17.704	2464	2468	2471	rBV4	86681	127552	1.41%	0.093%
83	18.004	2516	2519	2525	rVB4	68758	89612	0.99%	0.065%
84	18.104	2534	2536	2541	rVB5	56174	66613	0.74%	0.049%
85	18.163	2543	2546	2551	rBV2	80722	122415	1.35%	0.089%
86	18.222	2551	2556	2566	rBV	1328097	1933733	21.39%	1.408%
87	18.822	2654	2658	2661	rBV	79268	92967	1.03%	0.068%
88	19.116	2705	2708	2712	rBV5	45369	59758	0.66%	0.044%
89	19.163	2713	2716	2719	rVB3	56853	65235	0.72%	0.048%
90	19.357	2738	2749	2753	rBV3	306981	818737	9.05%	0.596%
91	19.451	2761	2765	2776	rBV	301362	515078	5.70%	0.375%
92	19.680	2799	2804	2817	rVB	5782714	8138561	90.01%	5.927%
93	20.533	2945	2949	2953	rBV4	220651	316448	3.50%	0.230%
94	20.674	2970	2973	2980	rBV4	175290	342175	3.78%	0.249%
95	21.133	3047	3051	3062	rBV	1394274	1922217	21.26%	1.400%
96	21.439	3098	3103	3108	rBV	4622050	6353495	70.26%	4.627%
97	22.080	3209	3212	3220	rVB4	449227	710704	7.86%	0.518%
98	23.139	3388	3392	3398	rBV2	711472	1116483	12.35%	0.813%
99	23.768	3492	3499	3515	rVB	4142499	9042304	100.00%	6.586%
100	23.927	3520	3526	3533	rVB2	3319021	6636928	73.40%	4.834%

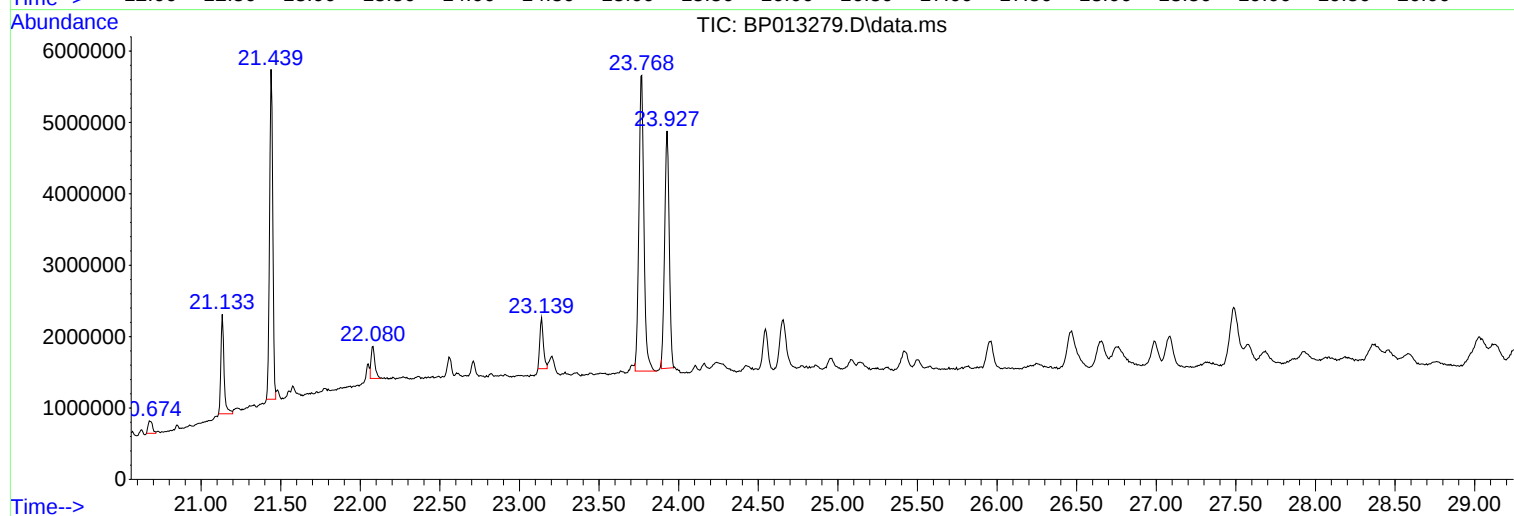
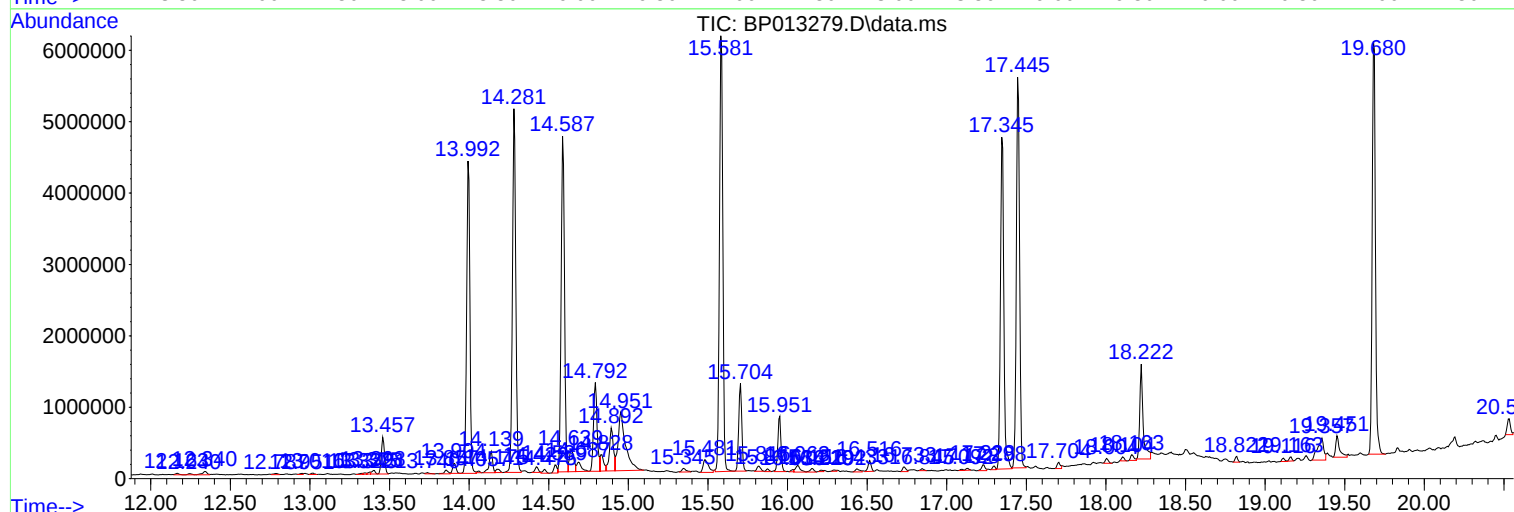
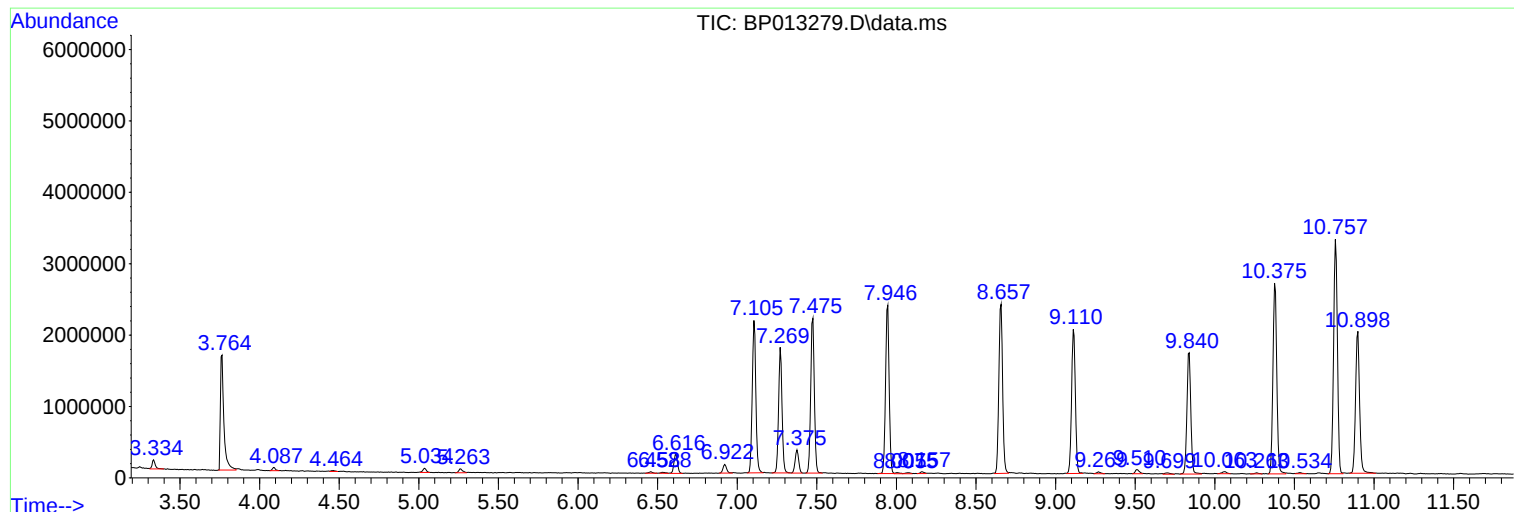
Sum of corrected areas: 137303148

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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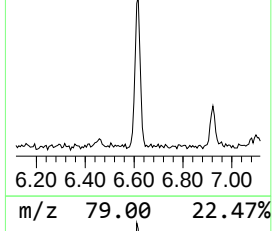
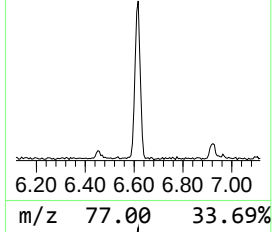
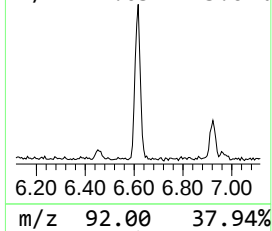
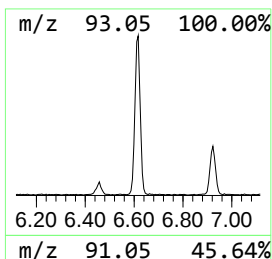
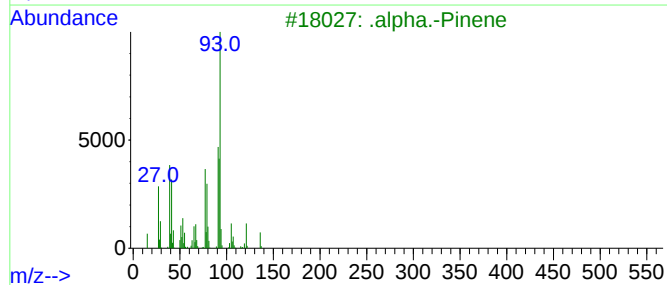
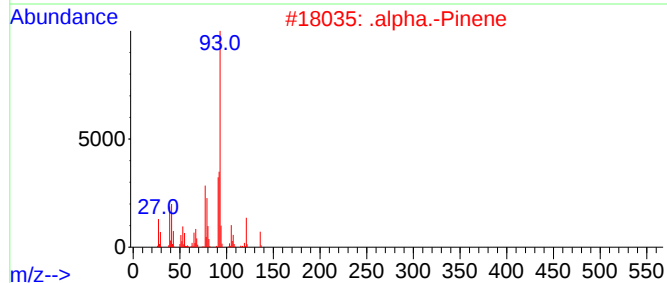
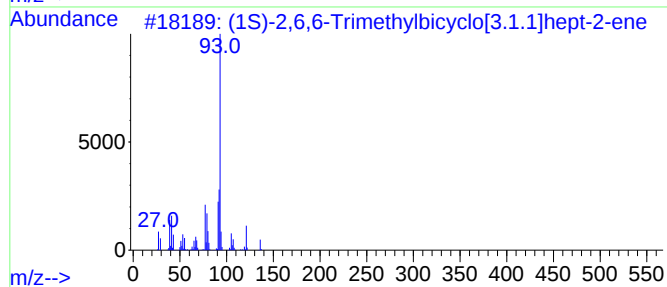
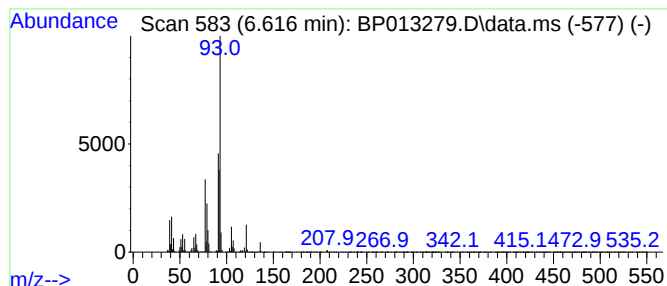
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (1S)-2,6,6-Trimethylbicyclo... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.616	2.22 ng/ul	415100	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1S)-2,6,6-Trimethylbicyclo[3.1...	136	C10H16	007785-26-4	95		
2	.alpha.-Pinene	136	C10H16	000080-56-8	95		
3	.alpha.-Pinene	136	C10H16	000080-56-8	94		
4	(1R)-2,6,6-Trimethylbicyclo[3.1...	136	C10H16	007785-70-8	91		
5	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	91		



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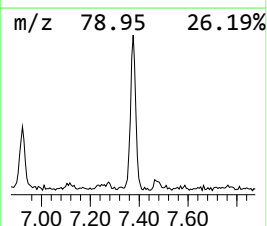
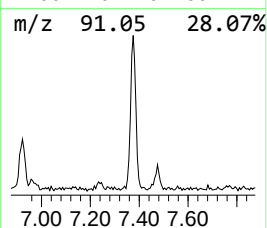
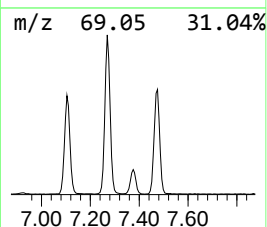
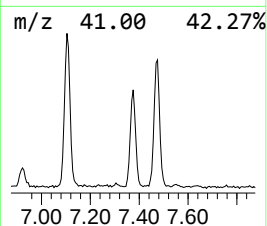
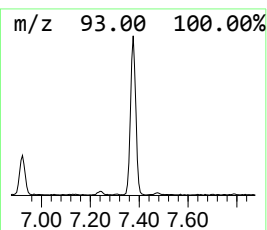
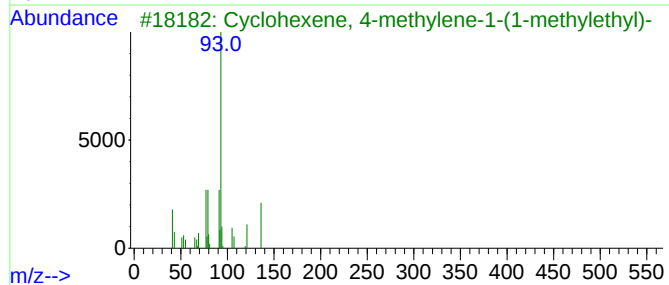
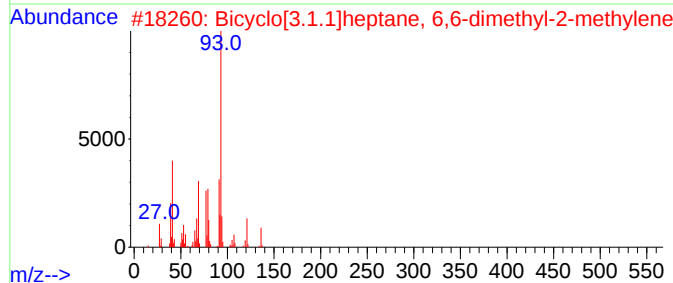
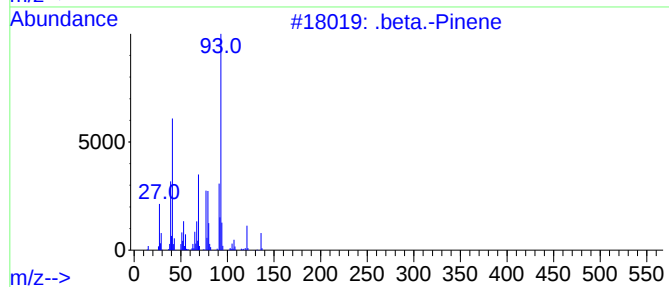
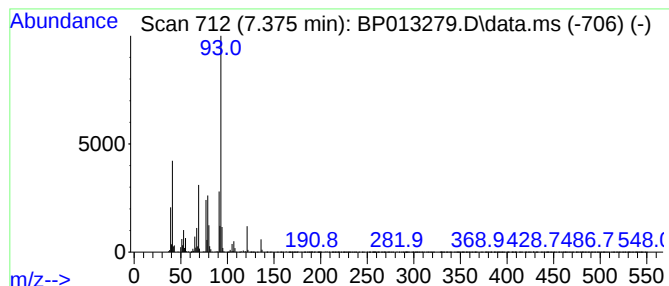
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 .beta.-Pinene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.375	2.64 ng/ul	493583	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	beta.-Pinene	136	C10H16	000127-91-3	96	
2	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	94		
3	Cyclohexene, 4-methylene-1-(1-me...	136	C10H16	000099-84-3	91		
4	.alpha.-Pinene	136	C10H16	000080-56-8	90		
5	.	beta.-Pinene	136	C10H16	000127-91-3	90	



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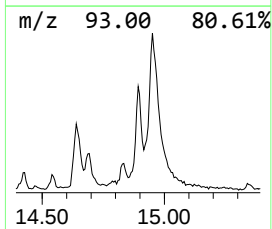
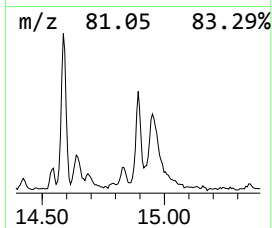
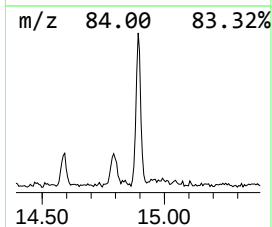
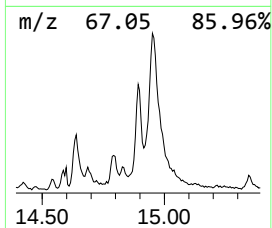
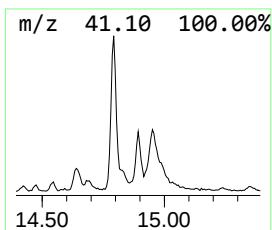
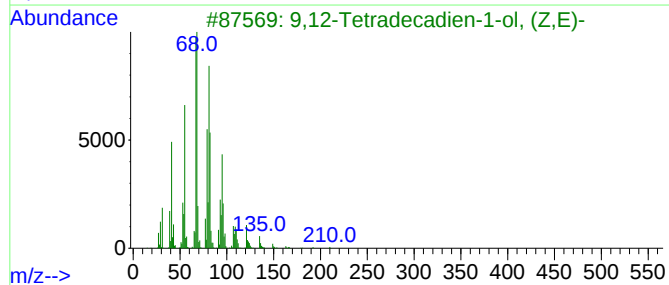
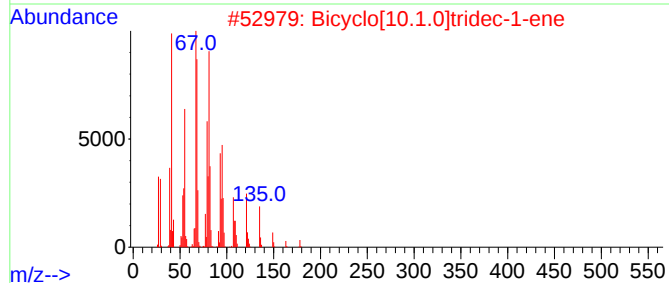
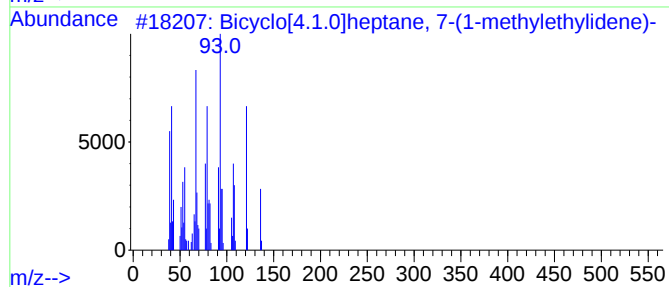
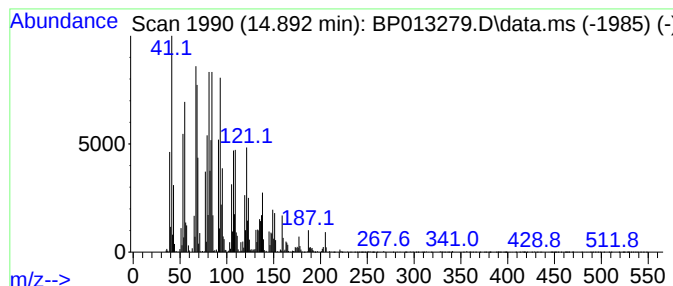
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Bicyclo[4.1.0]heptane, 7-(1... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.892	2.99 ng/ul	1028810	Acenaphthene-d10	14.586

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[4.1.0]heptane, 7-(1-meth...	136	C10H16	053282-47-6	60
2			Bicyclo[10.1.0]tridec-1-ene	178	C13H22	054766-91-5	46
3			9,12-Tetradecadien-1-ol, (Z,E)-	210	C14H26O	051937-00-9	46
4			o-Mentha-1(7),8-dien-3-ol	152	C10H16O	015358-81-3	42
5			Bicyclo[10.1.0]tridec-1-ene	178	C13H22	054766-91-5	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013279.D
Acq On : 23 Dec 2022 14:52
Operator : CG/JU
Sample : N6018-09
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYH9

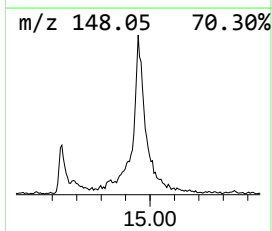
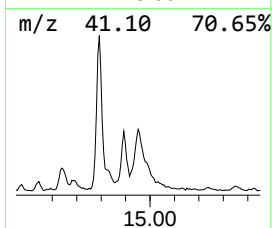
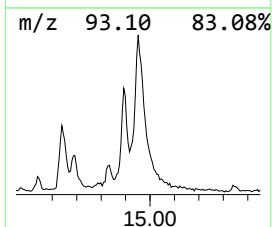
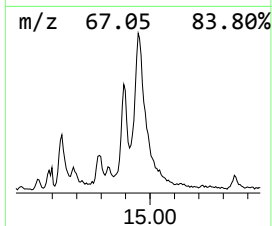
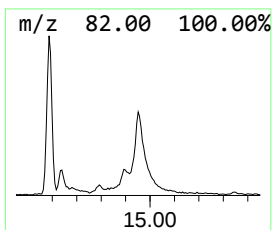
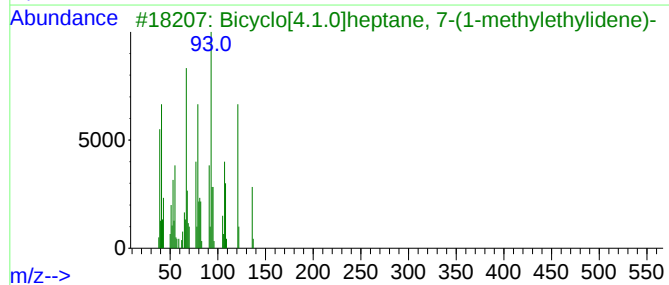
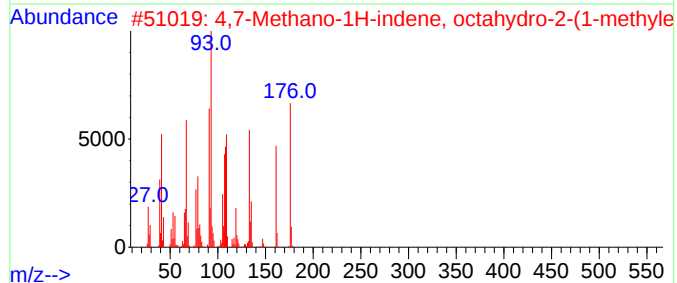
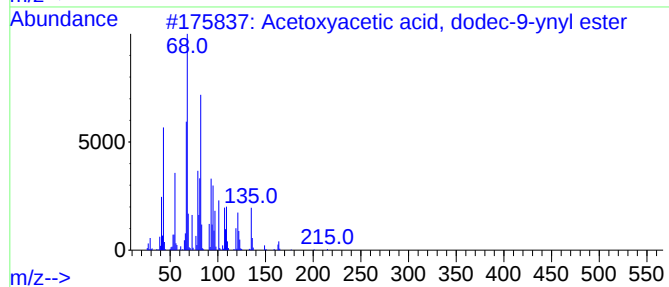
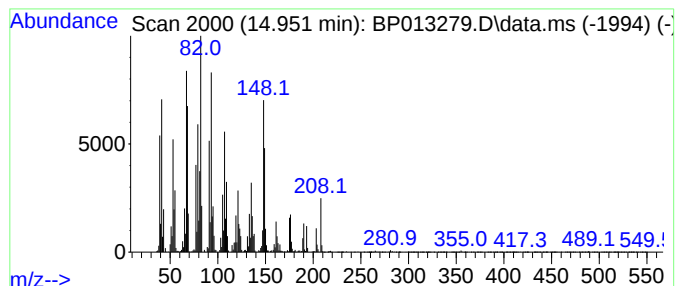
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.951	7.64 ng/ul	2628460	Acenaphthene-d10	14.586

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetoxyacetic acid, dodec-9-ynyl...	282	C16H26O4	1000308-31-9	38
2			4,7-Methano-1H-indene, octahydro...	176	C13H20	074793-54-7	38
3			Bicyclo[4.1.0]heptane, 7-(1-meth...	136	C10H16	053282-47-6	35
4			5-Chlorovaleric acid, dodec-9-yn...	300	C17H29ClO2	1000292-24-7	35
5			Camphene	136	C10H16	000079-92-5	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013279.D
Acq On : 23 Dec 2022 14:52
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Sample : N6018-09
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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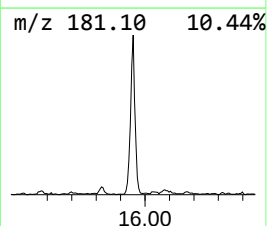
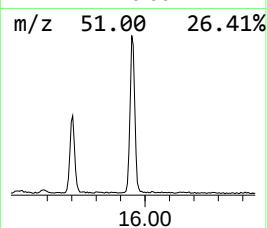
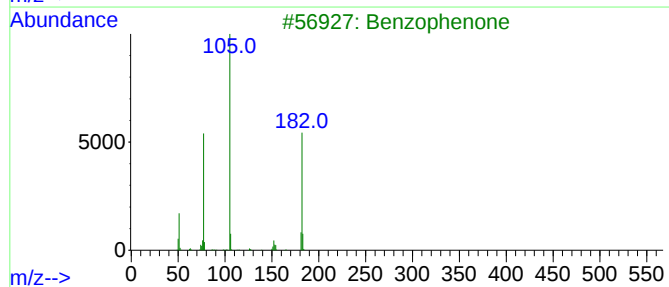
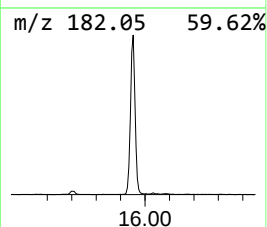
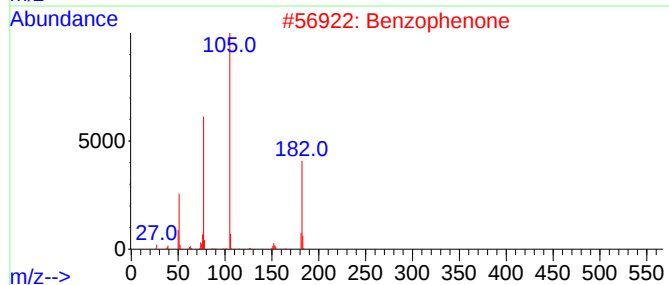
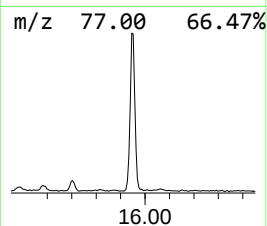
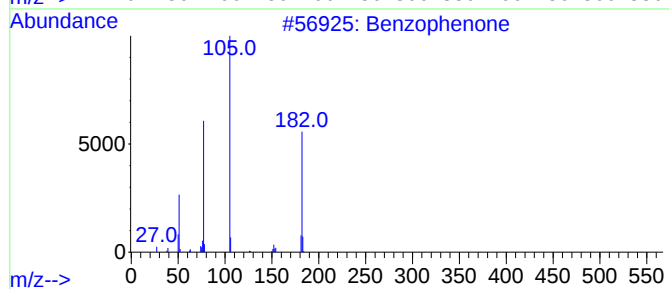
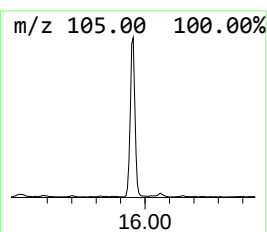
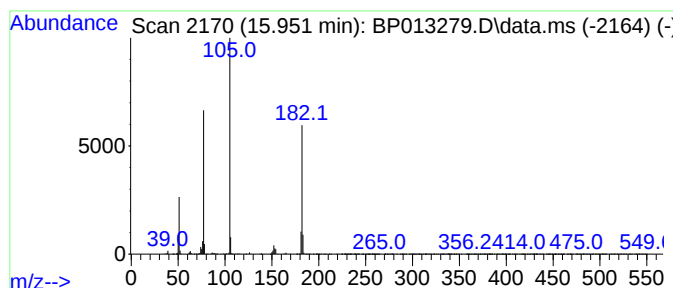
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.951	3.06 ng/ul	1052240	Acenaphthene-d10	14.586

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013279.D
Acq On : 23 Dec 2022 14:52
Operator : CG/JU
Sample : N6018-09
Misc :
ALS Vial : 47 Sample Multiplier: 1

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ClientSampleId :
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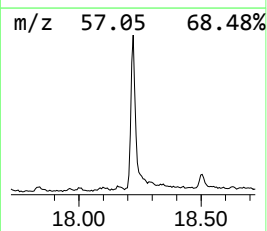
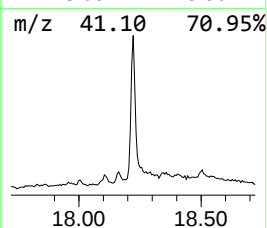
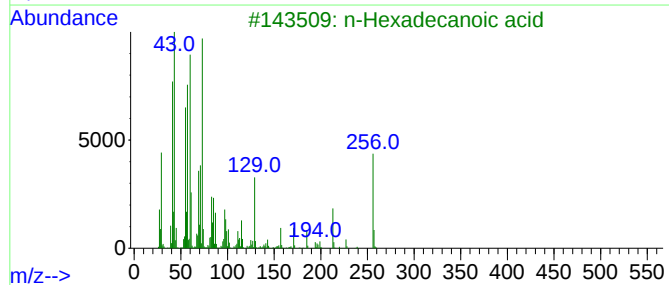
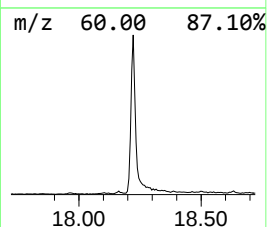
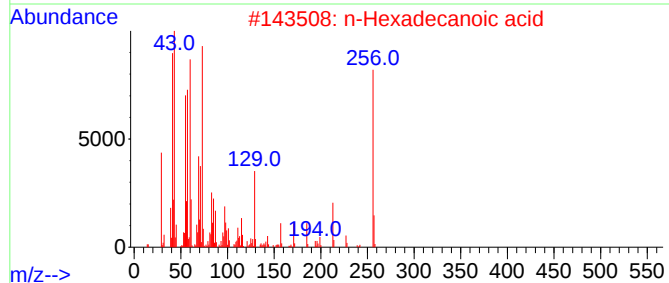
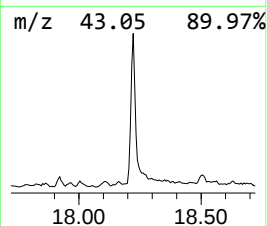
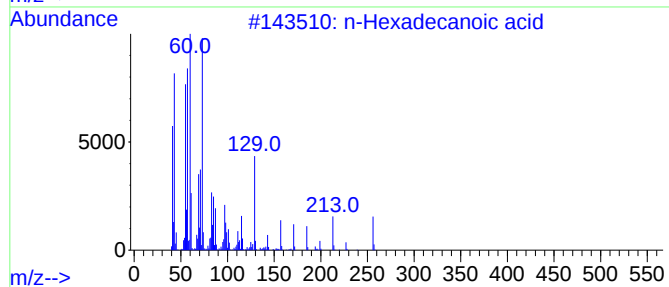
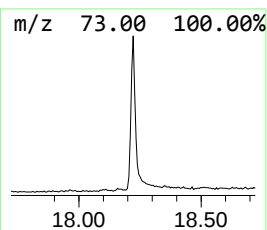
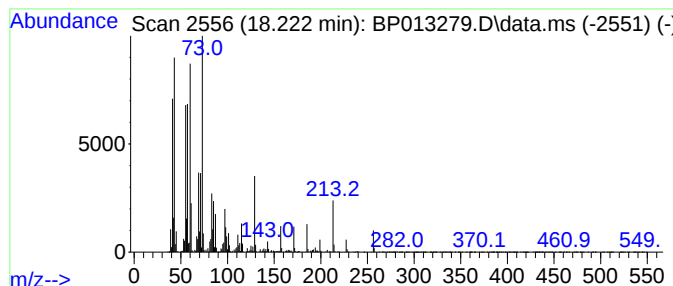
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.222	5.54 ng/ul	1933730	Phenanthrene-d10	17.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013279.D
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Sample : N6018-09
Misc :
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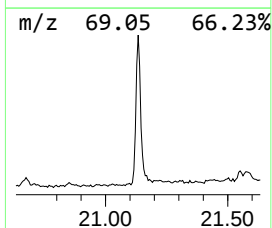
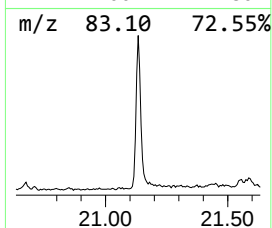
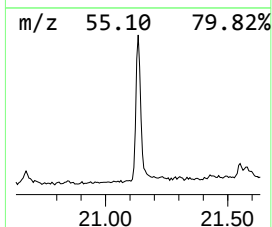
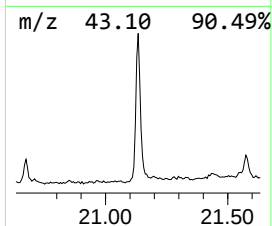
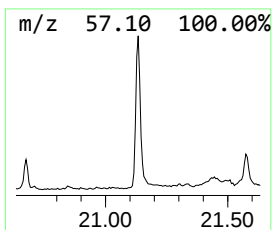
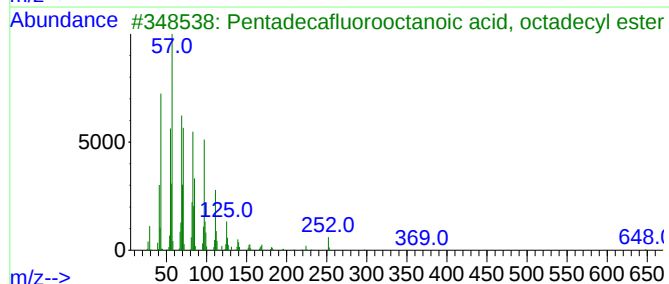
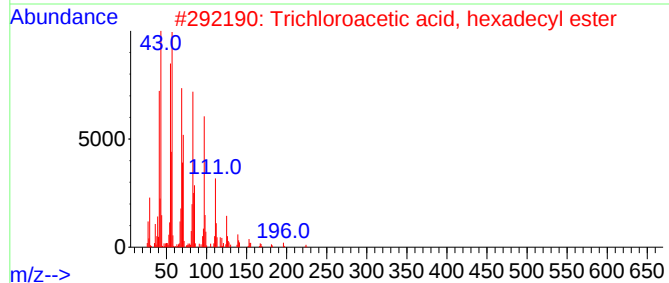
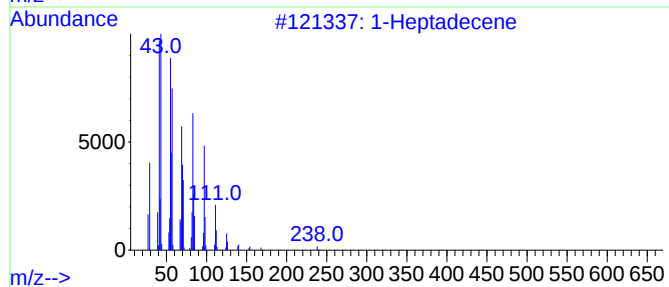
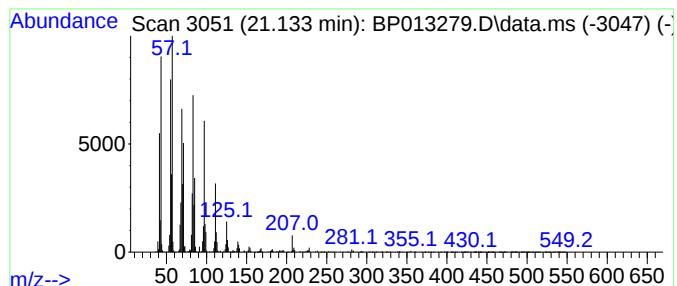
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.133	6.05 ng/ul	1922220	Chrysene-d12	21.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heptadecene	238	C17H34	006765-39-5	94
2			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
4			1-Hexacosene	364	C26H52	018835-33-1	94
5			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013279.D
Acq On : 23 Dec 2022 14:52
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Sample : N6018-09
Misc :
ALS Vial : 47 Sample Multiplier: 1

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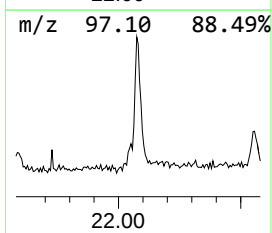
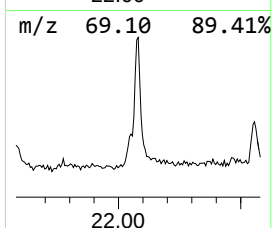
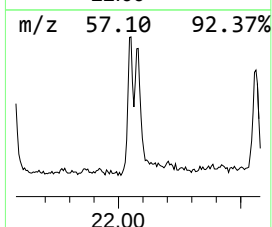
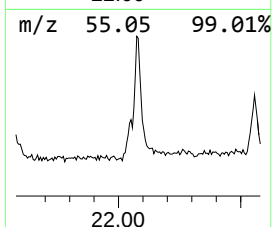
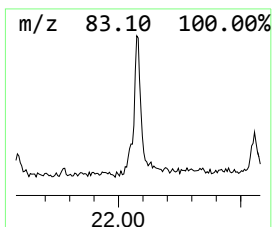
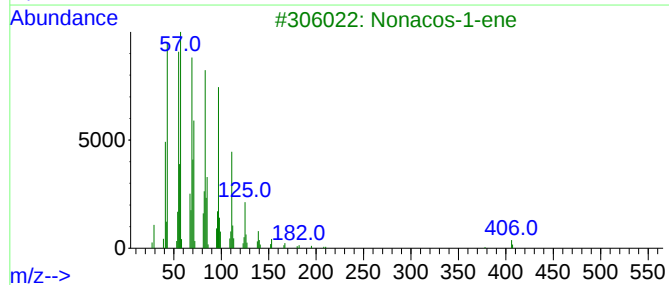
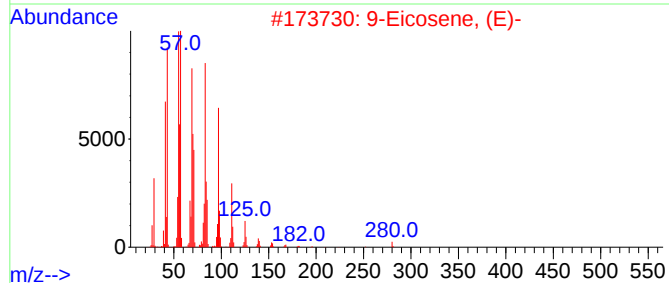
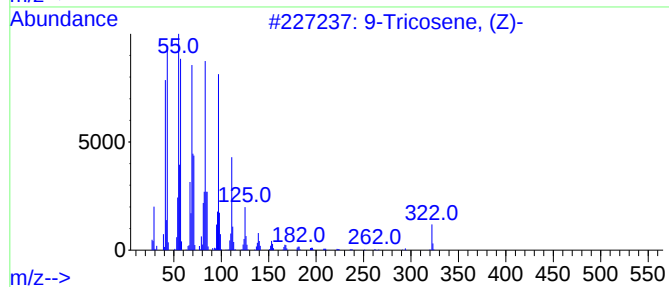
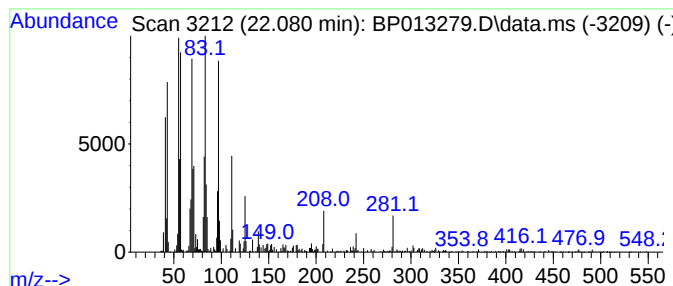
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.080	2.24 ng/ul	710704	Chrysene-d12	21.439

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Tricosene, (Z)-		322	C23H46	027519-02-4	87
2	9-Eicosene, (E)-		280	C20H40	074685-29-3	86
3	Nonacos-1-ene		406	C29H58	018835-35-3	70
4	5-Eicosene, (E)-		280	C20H40	074685-30-6	68
5	n-Tetracosanol-1		354	C24H50O	000506-51-4	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
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Sample : N6018-09
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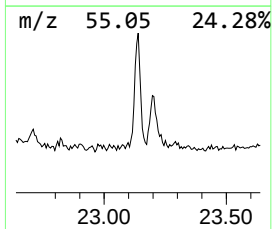
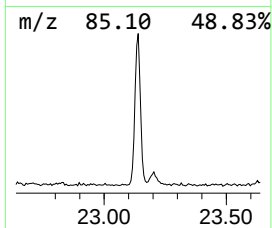
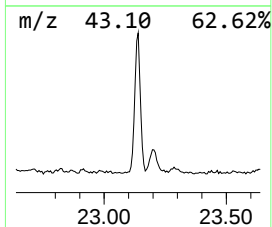
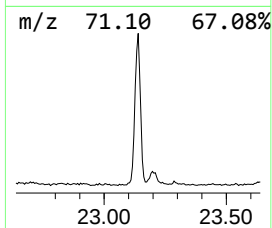
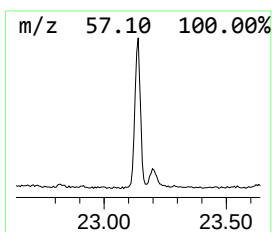
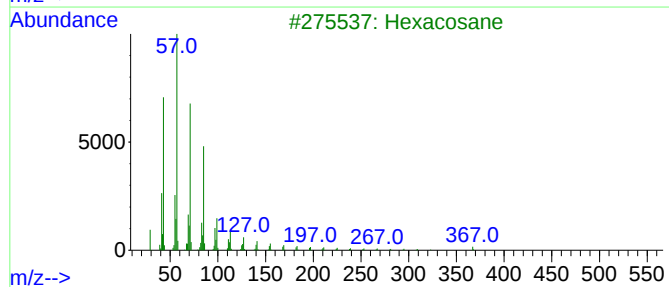
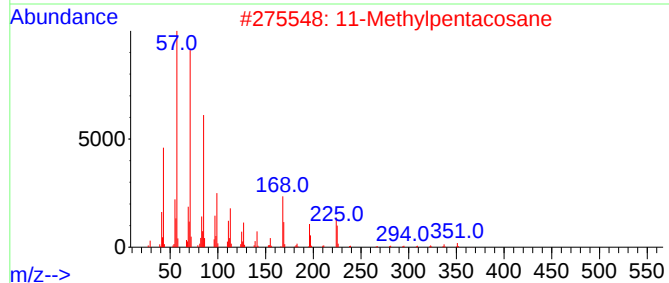
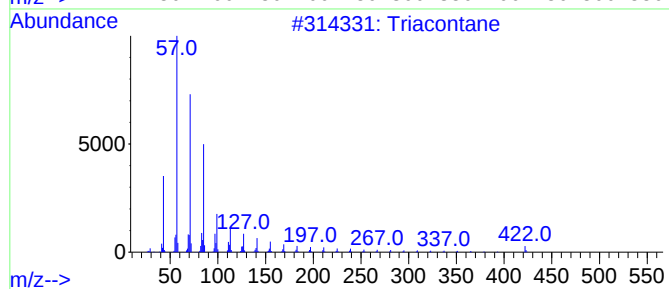
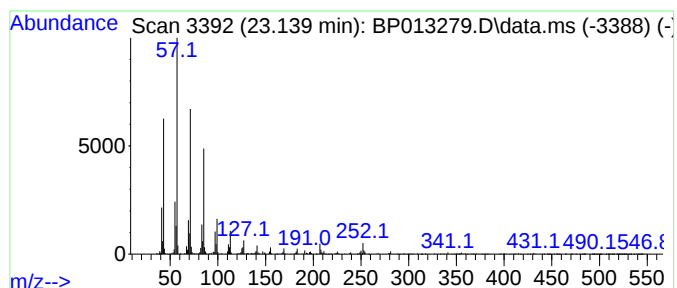
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.139	3.36 ng/ul	1116480	Perylene-d12	23.927

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Triacontane	422	C30H62	000638-68-6	97
2		11-Methylpentacosane	366	C26H54	015689-71-1	93
3		Hexacosane	366	C26H54	000630-01-3	93
4		Heneicosane	296	C21H44	000629-94-7	90
5		Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
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Operator : CG/JU
Sample : N6018-09
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYH9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.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
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.beta.-Pinene	7.375	2.6	ng/ul	493583	1	7.946	3732840	20.0
Cyclohexane, 1-...	13.457	2.1	ng/ul	717798	3	14.586	6878280	20.0
Bicyclo[4.1.0]h...	14.892	3.0	ng/ul	1028810	3	14.586	6878280	20.0
unknown-01	14.951	7.6	ng/ul	2628460	3	14.586	6878280	20.0
Benzophenone	15.951	3.1	ng/ul	1052240	3	14.586	6878280	20.0
n-Hexadecanoic ...	18.222	5.5	ng/ul	1933730	4	17.345	6983010	20.0
(DEL) Alkane: S...	21.133	6.0	ng/ul	1922220	5	21.439	6353500	20.0
(DEL) Alkane: S...	22.080	2.2	ng/ul	710704	5	21.439	6353500	20.0
(DEL) Alkane: S...	23.139	3.4	ng/ul	1116480	6	23.927	6636930	20.0