

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
 Data File : BP013296.D
 Acq On : 24 Dec 2022 13:00
 Operator : CG/JU
 Sample : N6016-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBYC0

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: BP013296.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	33	rVB	114038	162700	1.88%	0.132%
2	3.775	99	100	109	rVB	91001	112613	1.30%	0.091%
3	3.864	111	115	122	rVB	130208	177800	2.05%	0.144%
4	3.987	132	136	143	rVB	121339	184183	2.13%	0.149%
5	4.087	149	153	162	rVB	54105	79575	0.92%	0.065%
6	4.464	215	217	225	rVB3	27472	38750	0.45%	0.031%
7	4.652	246	249	253	rVB	47470	56882	0.66%	0.046%
8	4.699	253	257	265	rVB	143110	198710	2.29%	0.161%
9	5.017	306	311	320	rBV3	394966	780703	9.01%	0.633%
10	5.117	325	328	333	rVB5	20650	31489	0.36%	0.026%
11	5.499	391	393	399	rVB6	13298	17243	0.20%	0.014%
12	5.622	411	414	419	rVB	16298	21350	0.25%	0.017%
13	6.540	566	570	575	rBV	43655	69866	0.81%	0.057%
14	6.611	577	582	590	rVB	45971	76498	0.88%	0.062%
15	6.916	630	634	638	rBV7	7726	13639	0.16%	0.011%
16	7.105	656	666	680	rVB	1759619	2901932	33.49%	2.353%
17	7.269	687	694	702	rBV	1451482	2255528	26.03%	1.829%
18	7.369	709	711	716	rVB6	9762	12844	0.15%	0.010%
19	7.469	721	728	738	rVB	1873872	2901389	33.49%	2.353%
20	7.546	738	741	746	rBV	28553	41393	0.48%	0.034%
21	7.752	772	776	782	rVB2	12023	23091	0.27%	0.019%
22	7.940	801	808	815	rBV	2618761	4047214	46.71%	3.282%
23	7.999	816	818	824	rVB6	16108	21779	0.25%	0.018%
24	8.652	922	929	938	rBV	1805245	2855418	32.96%	2.315%
25	8.769	943	949	956	rVB	236909	383469	4.43%	0.311%
26	9.110	1000	1007	1020	rBV	1737282	2871358	33.14%	2.328%
27	9.269	1029	1034	1042	rVB10	15447	27836	0.32%	0.023%
28	9.505	1069	1074	1083	rVB	75849	125283	1.45%	0.102%
29	9.834	1122	1130	1144	rBV	1451131	2421114	27.94%	1.963%
30	9.987	1151	1156	1161	rBV2	20646	38849	0.45%	0.032%
31	10.057	1163	1168	1173	rVB8	12130	23519	0.27%	0.019%
32	10.193	1186	1191	1196	rBV2	22213	38833	0.45%	0.031%
33	10.375	1215	1222	1235	rBV	2192094	3681409	42.49%	2.985%
34	10.757	1278	1287	1303	rBV	3544886	6116196	70.59%	4.960%
35	10.893	1303	1310	1328	rVV	1339284	2458408	28.37%	1.994%
36	11.287	1371	1377	1383	rBV7	21510	40854	0.47%	0.033%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
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37	11.552	1415	1422	1431	rVB4	19991	47559	0.55%	0.039%
38	12.004	1495	1499	1503	rVB6	12532	21240	0.25%	0.017%
39	12.116	1515	1518	1522	rBV6	7184	14363	0.17%	0.012%
40	12.163	1522	1526	1531	rVB	30492	43343	0.50%	0.035%
41	12.246	1537	1540	1544	rVB4	14968	20176	0.23%	0.016%
42	12.334	1551	1555	1562	rVB	38832	66249	0.76%	0.054%
43	13.110	1682	1687	1692	rBV2	32682	51626	0.60%	0.042%
44	13.187	1694	1700	1703	rBV8	9118	16422	0.19%	0.013%
45	13.398	1724	1736	1740	rBV	72332	116283	1.34%	0.094%
46	13.551	1758	1762	1768	rVB7	9726	21297	0.25%	0.017%
47	13.904	1817	1822	1827	rBV8	27691	40948	0.47%	0.033%
48	13.993	1829	1837	1846	rBV	3776524	5186593	59.86%	4.206%
49	14.104	1853	1856	1860	rVB5	15469	18699	0.22%	0.015%
50	14.228	1874	1877	1879	rBV4	11707	16636	0.19%	0.013%
51	14.281	1879	1886	1904	rVB	4181468	6221916	71.81%	5.045%
52	14.469	1914	1918	1923	rVB2	43306	61599	0.71%	0.050%
53	14.587	1931	1938	1945	rVV	5385646	7705629	88.93%	6.249%
54	14.645	1945	1948	1952	rVB6	13524	17610	0.20%	0.014%
55	14.787	1966	1972	1990	rBV	870371	1503304	17.35%	1.219%
56	14.940	1994	1998	2002	rVV3	43604	75142	0.87%	0.061%
57	14.998	2002	2008	2018	rVB4	97825	181378	2.09%	0.147%
58	15.175	2032	2038	2044	rBV	141643	212429	2.45%	0.172%
59	15.245	2047	2050	2054	rVB6	8919	14339	0.17%	0.012%
60	15.363	2065	2070	2074	rBV6	14344	31796	0.37%	0.026%
61	15.422	2078	2080	2083	rVB4	18894	18253	0.21%	0.015%
62	15.581	2100	2107	2113	rBV	5207024	7442269	85.89%	6.035%
63	15.634	2113	2116	2121	rVV4	35023	56736	0.65%	0.046%
64	15.698	2121	2127	2144	rVV	870686	1258486	14.52%	1.021%
65	15.816	2144	2147	2154	rVB4	29114	49279	0.57%	0.040%
66	15.910	2160	2163	2165	rVV3	9436	14318	0.17%	0.012%
67	15.945	2165	2169	2181	rVB	85208	136094	1.57%	0.110%
68	16.287	2224	2227	2234	rVB9	20632	41872	0.48%	0.034%
69	16.728	2298	2302	2310	rVB6	31610	40137	0.46%	0.033%
70	16.875	2324	2327	2332	rVB3	24489	30496	0.35%	0.025%
71	16.998	2343	2348	2354	rBV10	20411	38313	0.44%	0.031%
72	17.087	2359	2363	2367	rBV5	21357	34646	0.40%	0.028%
73	17.128	2367	2370	2375	rVB4	16049	26014	0.30%	0.021%
74	17.228	2383	2387	2394	rBV9	27030	40756	0.47%	0.033%
75	17.292	2395	2398	2399	rBV3	16710	16073	0.19%	0.013%
76	17.345	2400	2407	2412	rBV	5238918	7450359	85.99%	6.042%

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77	17.445	2418	2424	2432	rBV	4343470	6304797	72.77%	5.113%
78	17.534	2436	2439	2448	rVB8	48618	84285	0.97%	0.068%
79	17.616	2450	2453	2460	rVB6	25495	44365	0.51%	0.036%
80	17.998	2515	2518	2523	rBV7	45253	59175	0.68%	0.048%
81	18.098	2531	2535	2541	rBV6	35486	64689	0.75%	0.052%
82	18.157	2542	2545	2548	rBV2	42654	50191	0.58%	0.041%
83	18.216	2550	2555	2565	rBV	628773	917632	10.59%	0.744%
84	18.410	2584	2588	2593	rVB2	80539	110581	1.28%	0.090%
85	18.551	2608	2612	2617	rBV2	83191	122215	1.41%	0.099%
86	18.628	2622	2625	2630	rVB2	67783	85335	0.98%	0.069%
87	18.781	2648	2651	2657	rVB	137426	166636	1.92%	0.135%
88	19.251	2729	2731	2737	rVB4	92610	128684	1.49%	0.104%
89	19.351	2745	2748	2753	rVB	219636	279898	3.23%	0.227%
90	19.445	2761	2764	2771	rBV2	169739	236715	2.73%	0.192%
91	19.675	2798	2803	2812	rBV	5396714	7536741	86.99%	6.112%
92	19.828	2825	2829	2836	rBV	528783	616651	7.12%	0.500%
93	21.127	3046	3050	3056	rBV	1231290	1895489	21.88%	1.537%
94	21.433	3097	3102	3107	rBV	5622001	7520205	86.79%	6.098%
95	22.074	3207	3211	3220	rVB2	825617	1375898	15.88%	1.116%
96	23.127	3387	3390	3396	rBV2	645120	1047275	12.09%	0.849%
97	23.192	3398	3401	3413	rVB2	920128	1960860	22.63%	1.590%
98	23.757	3491	3497	3511	rVB2	4046528	8664404	100.00%	7.026%
99	23.921	3518	3525	3537	rVB2	3922192	8416832	97.14%	6.825%
100	24.533	3625	3629	3638	rVB	938157	1918948	22.15%	1.556%

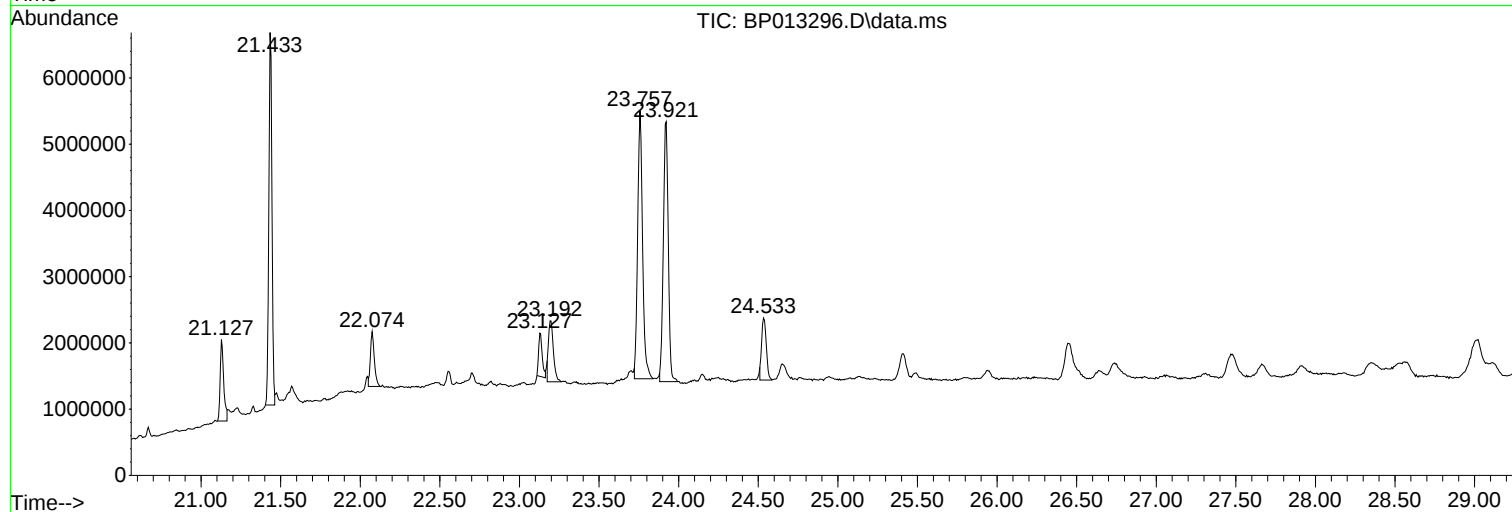
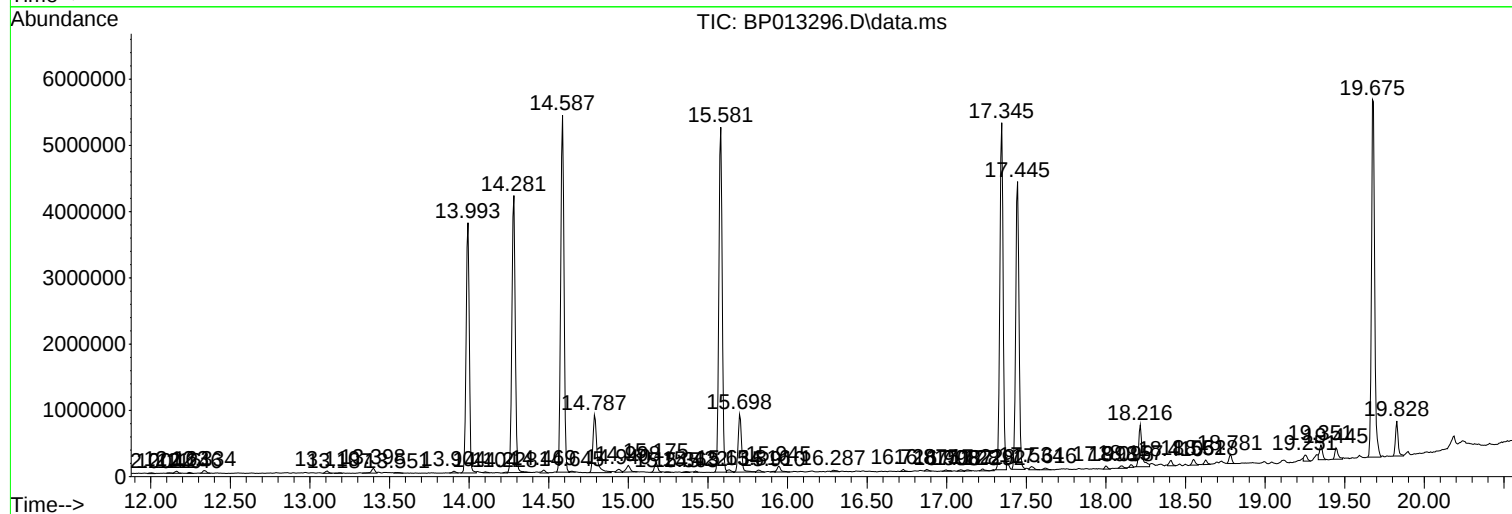
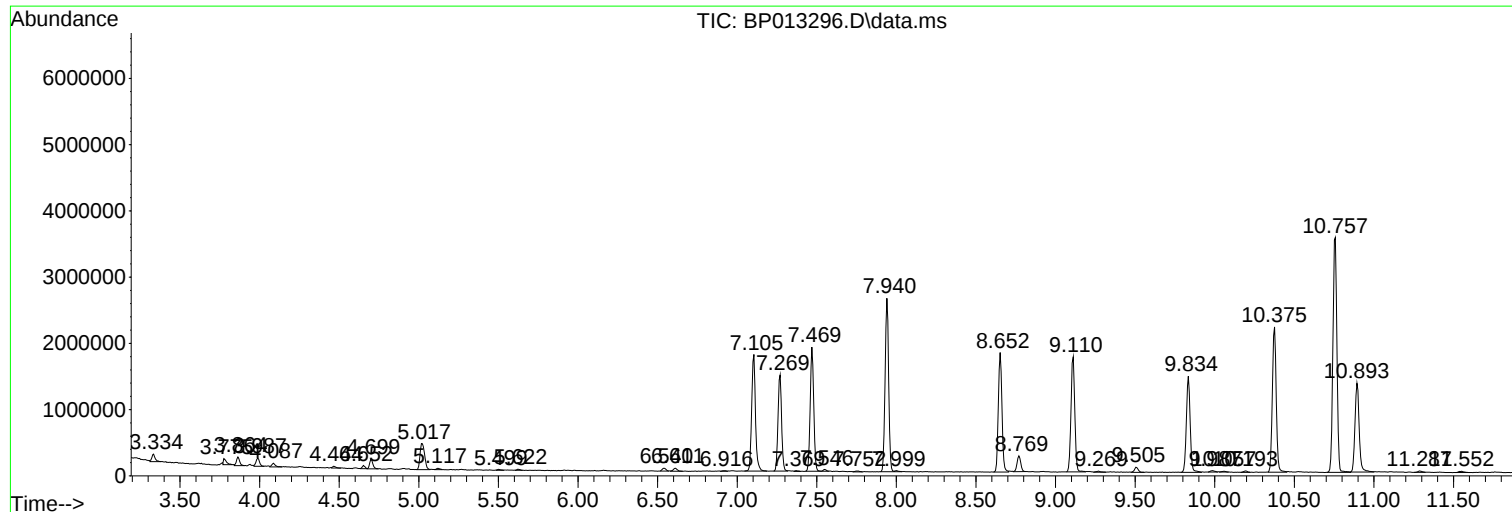
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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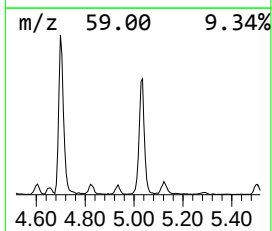
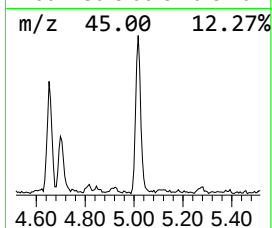
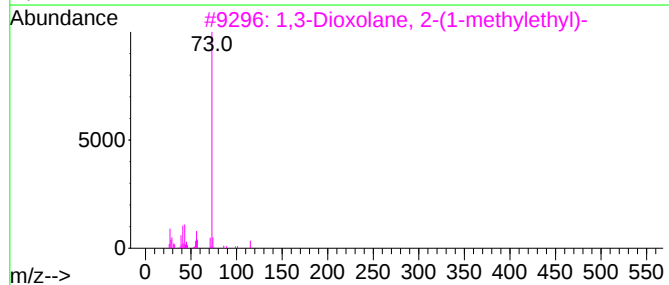
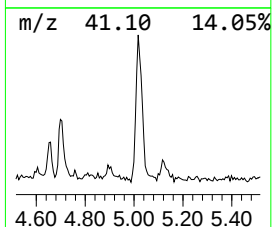
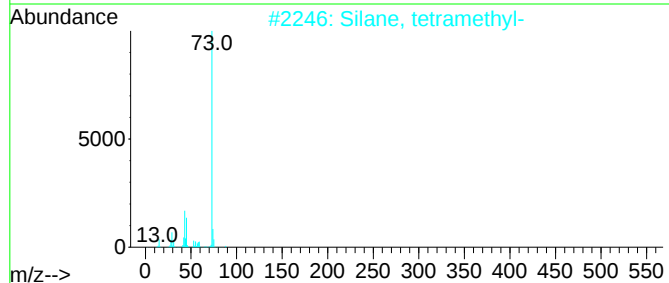
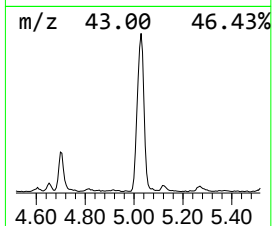
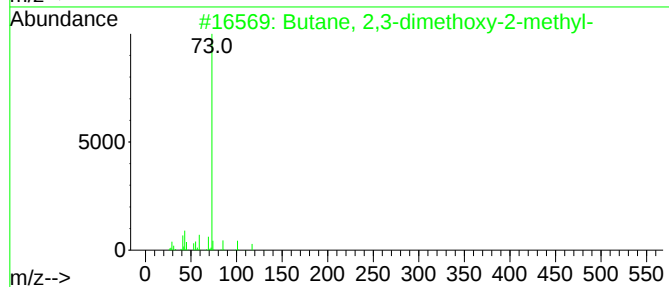
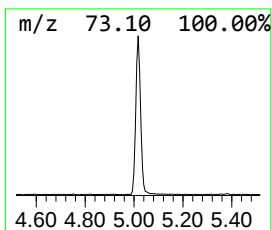
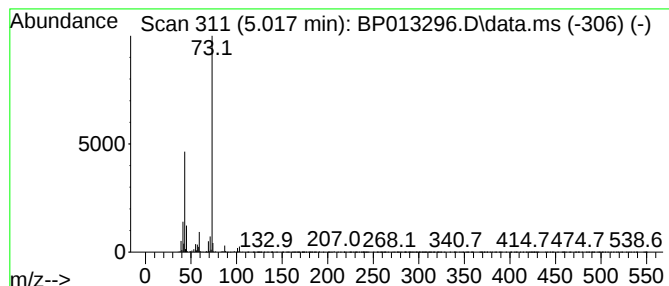
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.017	3.86 ng/ul	780703	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethoxy-2-methyl-	132	C7H16O2	074421-00-4	36		
2	Silane, tetramethyl-	88	C4H12Si	000075-76-3	32		
3	1,3-Dioxolane, 2-(1-methylethyl)-	116	C6H12O2	000822-83-3	23		
4	2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	23		
5	1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	17		



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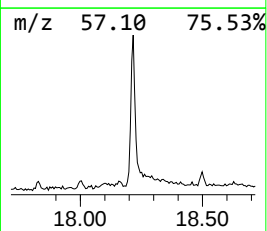
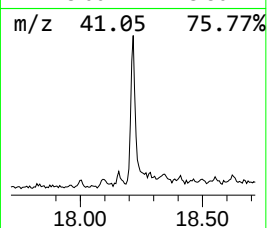
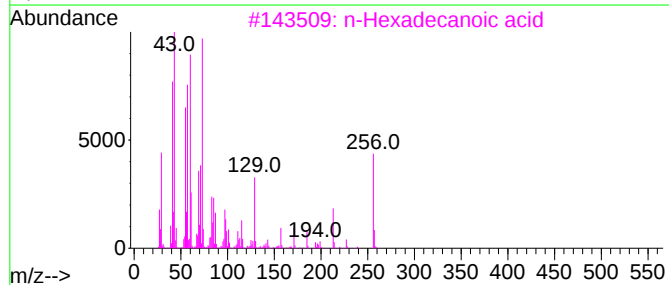
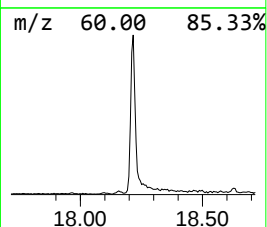
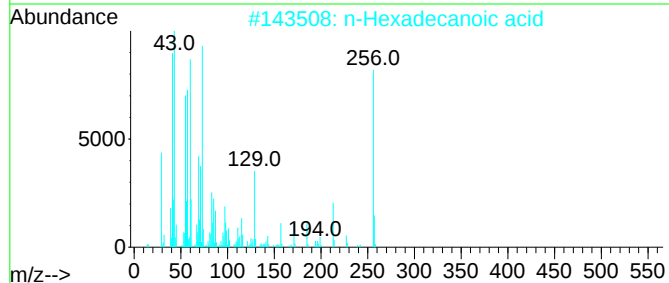
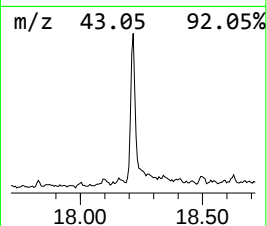
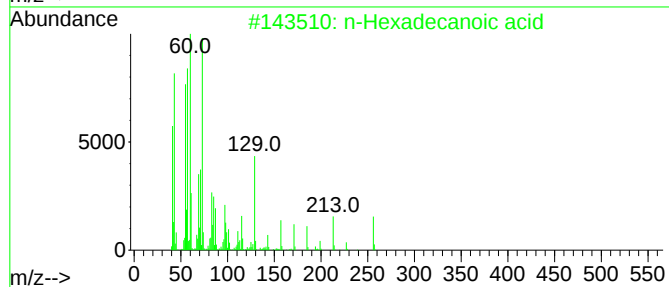
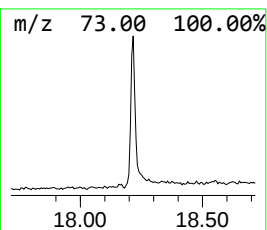
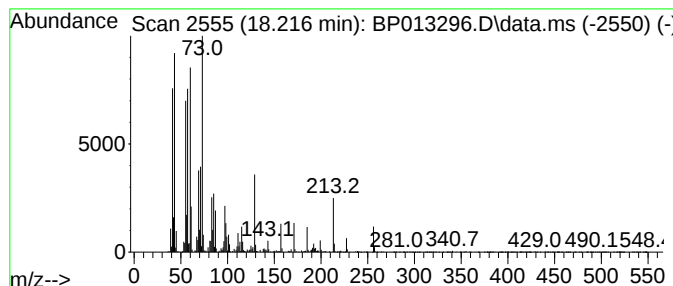
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	2.46 ng/ul	917632	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	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	87



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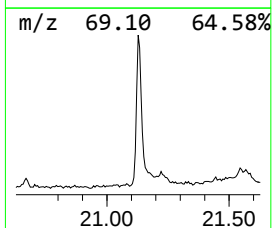
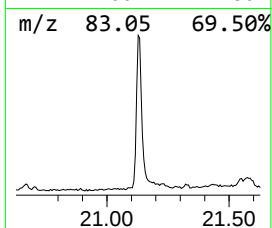
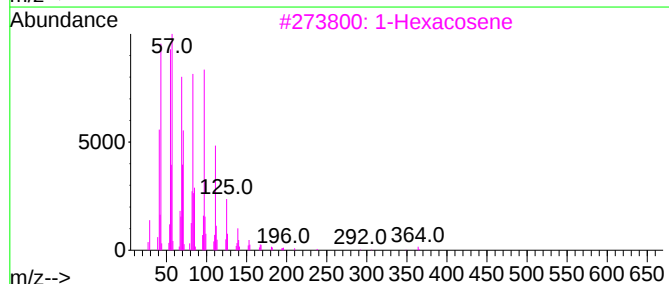
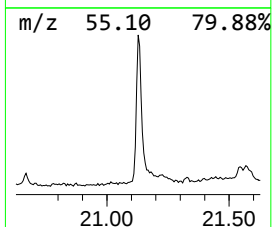
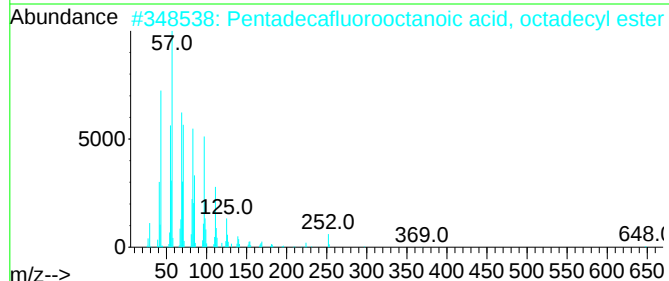
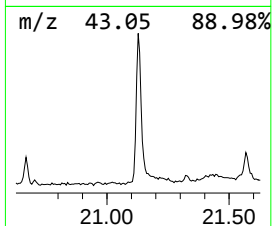
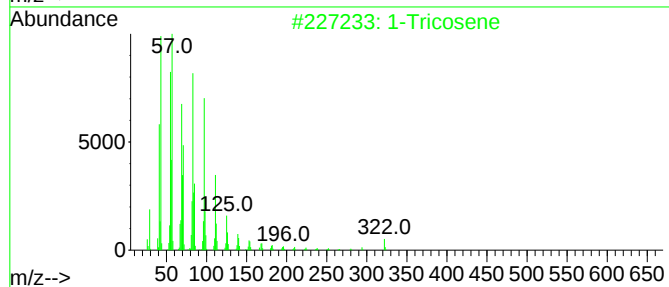
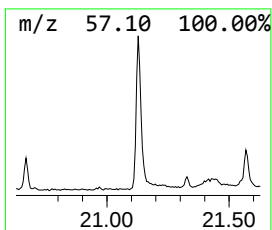
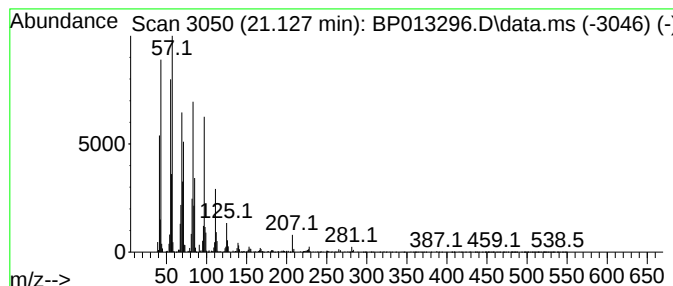
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ClientSampleId :
DBYC0

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 3 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.128	5.04 ng/ul	1895490	Chrysene-d12	21.433	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tricosene	322	C23H46	018835-32-0	95
2	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
3	1-Hexacosene	364	C26H52	018835-33-1	94
4	9-Tricosene, (Z)-	322	C23H46	027519-02-4	93
5	Pentacos-1-ene	350	C25H50	016980-85-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013296.D
Acq On : 24 Dec 2022 13:00
Operator : CG/JU
Sample : N6016-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYC0

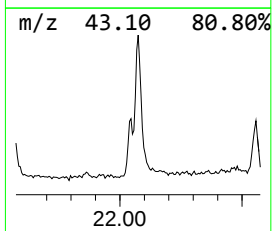
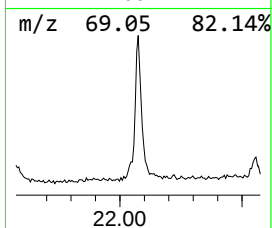
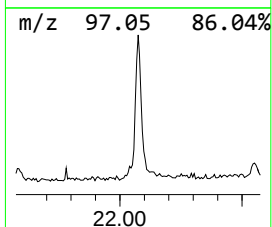
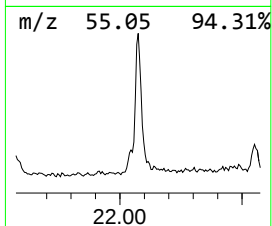
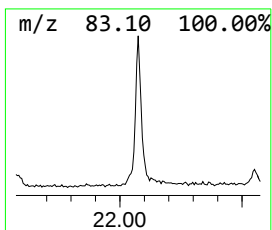
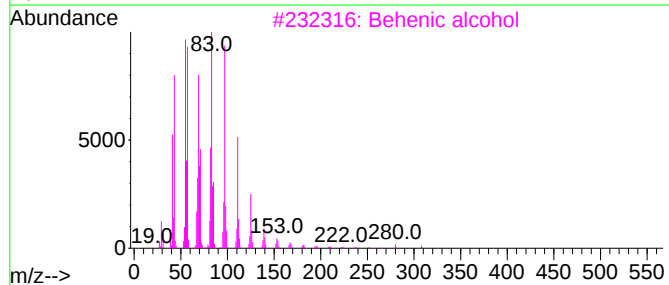
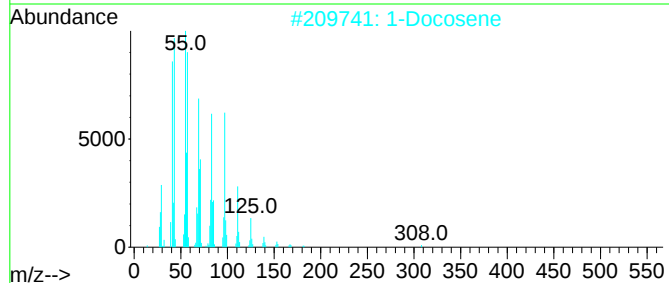
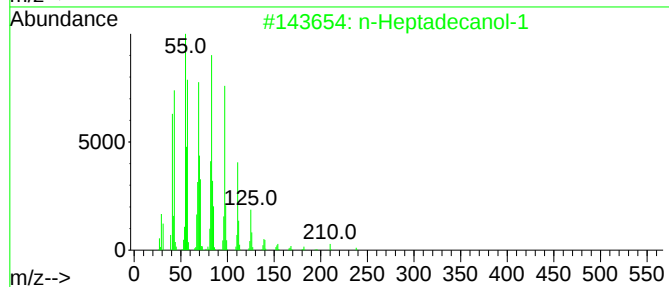
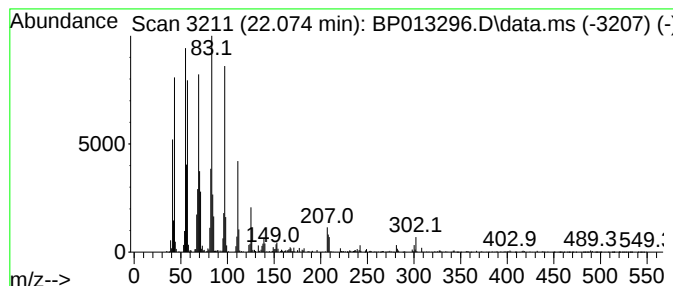
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Heptadecanol-1 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.075	3.66 ng/ul	1375900	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Heptadecanol-1	256	C17H36O	001454-85-9	94
2			1-Docosene	308	C22H44	001599-67-3	91
3			Behenic alcohol	326	C22H46O	000661-19-8	91
4			1-Hexadecanol	242	C16H34O	036653-82-4	91
5			1-Heneicosyl formate	340	C22H44O2	077899-03-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013296.D
Acq On : 24 Dec 2022 13:00
Operator : CG/JU
Sample : N6016-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

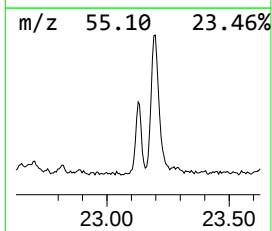
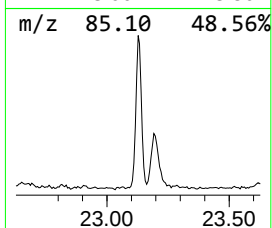
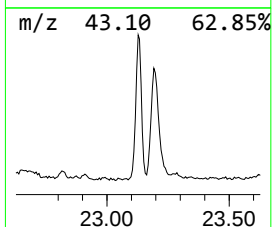
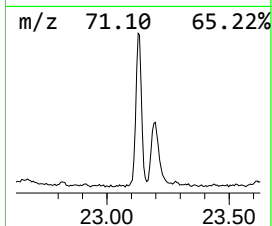
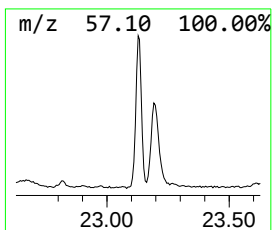
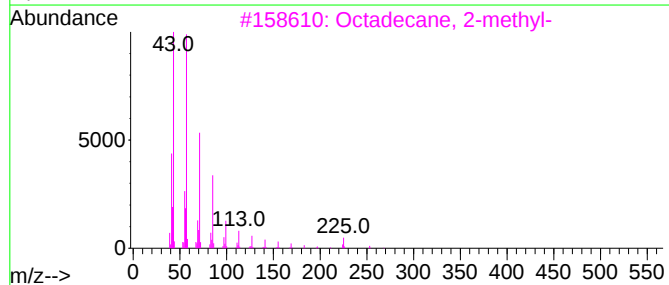
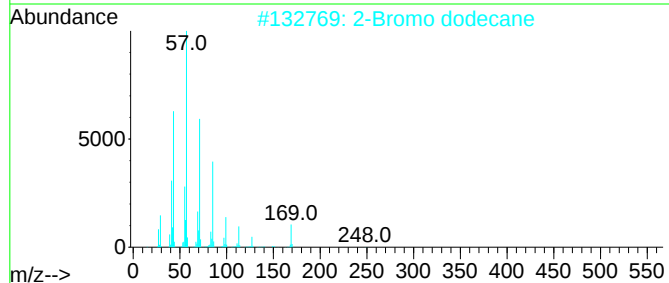
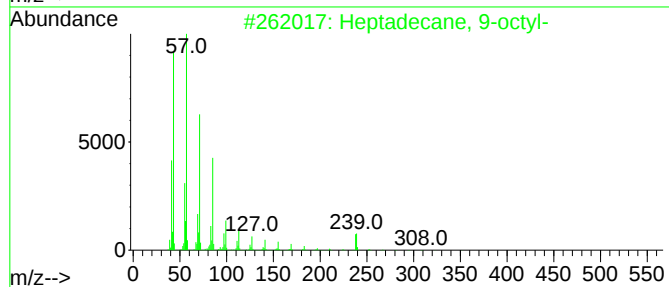
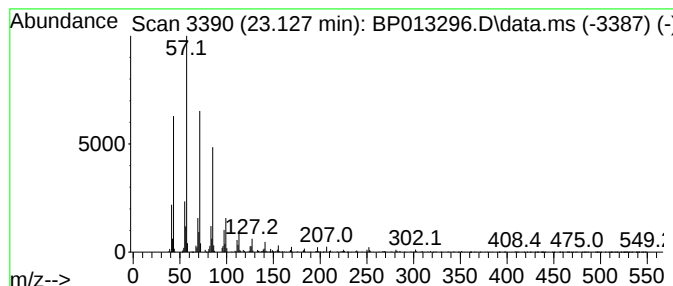
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.127	2.49 ng/ul	1047280	Perylene-d12	23.921	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
2	2-Bromo dodecane	248	C12H25Br	013187-99-0	93
3	Octadecane, 2-methyl-	268	C19H40	001560-88-9	93
4	Eicosane, 2-methyl-	296	C21H44	001560-84-5	91
5	Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013296.D
Acq On : 24 Dec 2022 13:00
Operator : CG/JU
Sample : N6016-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYC0

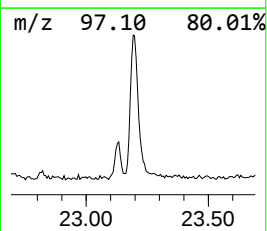
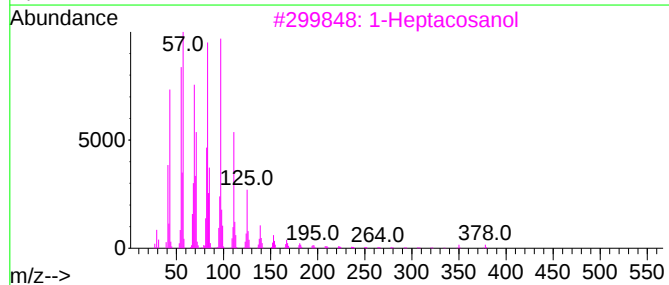
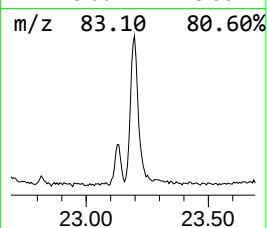
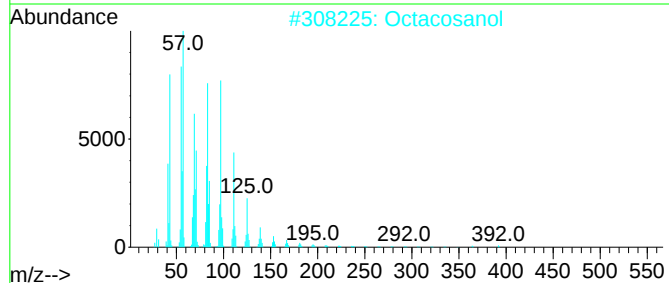
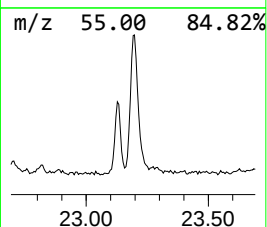
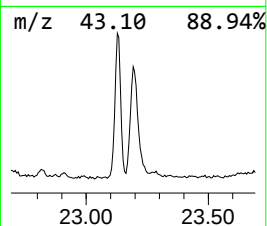
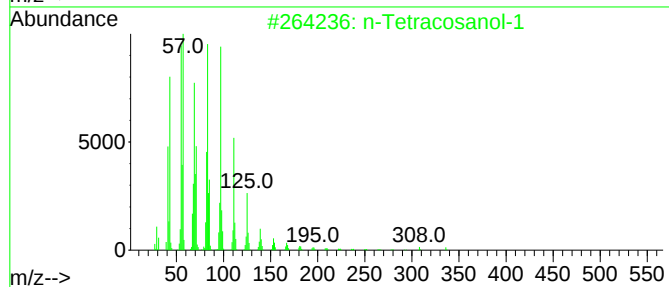
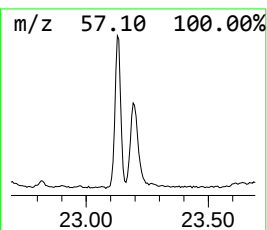
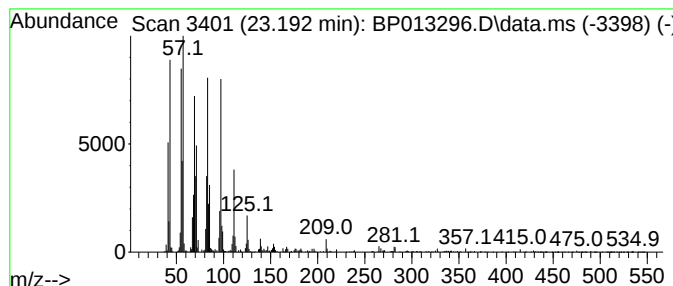
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Tetracosanol-1 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.192	4.66 ng/ul	1960860	Perylene-d12	23.921

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
2			Octacosanol	410	C28H58O	000557-61-9	91
3			1-Heptacosanol	396	C27H56O	002004-39-9	91
4			1-Tetracosene	336	C24H48	010192-32-2	91
5			1-Docosene	308	C22H44	001599-67-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013296.D
Acq On : 24 Dec 2022 13:00
Operator : CG/JU
Sample : N6016-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

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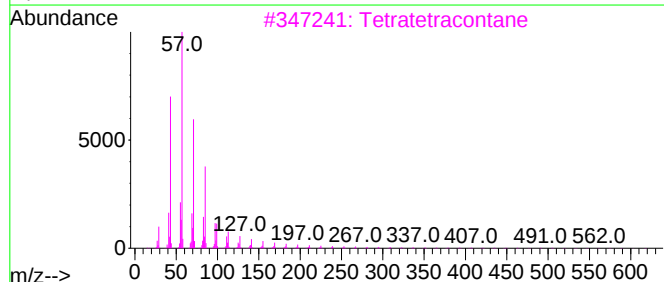
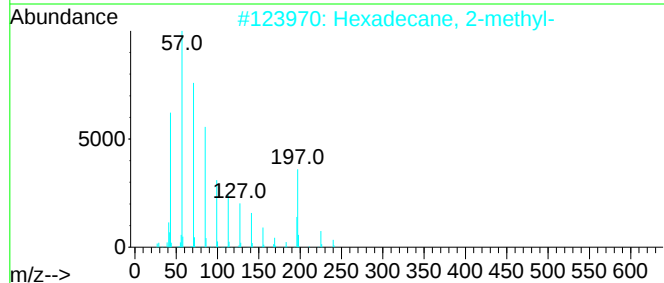
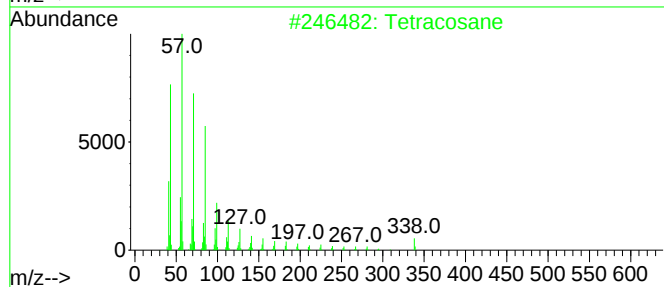
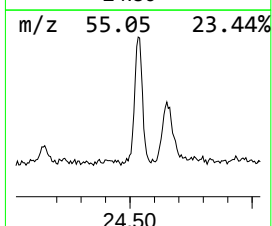
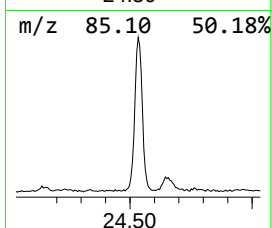
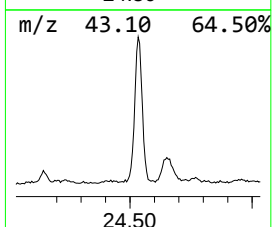
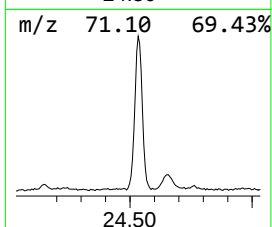
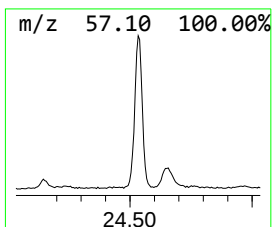
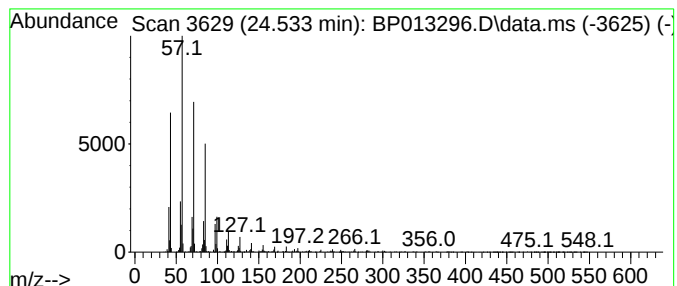
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.533	4.56 ng/ul	1918950	Perylene-d12	23.921

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	96
2			Hexadecane, 2-methyl-	240	C17H36	001560-92-5	91
3			Tetratetracontane	619	C44H90	007098-22-8	91
4			Heneicosane	296	C21H44	000629-94-7	90
5			Octacosane, 2-methyl-	408	C29H60	001560-98-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013296.D
Acq On : 24 Dec 2022 13:00
Operator : CG/JU
Sample : N6016-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYC0

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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n-Hexadecanoic ...	18.216	2.5	ng/ul	917632	4	17.345	7450360	20.0
(DEL) Alkane: S...	21.128	5.0	ng/ul	1895490	5	21.433	7520210	20.0
n-Heptadecanol-1	22.075	3.7	ng/ul	1375900	5	21.433	7520210	20.0
(DEL) Alkane: S...	23.127	2.5	ng/ul	1047280	6	23.921	8416830	20.0
n-Tetracosanol-1	23.192	4.7	ng/ul	1960860	6	23.921	8416830	20.0
(DEL) Alkane: S...	24.533	4.6	ng/ul	1918950	6	23.921	8416830	20.0