

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
 Data File : BP013310.D
 Acq On : 24 Dec 2022 21:28
 Operator : CG/JU
 Sample : N6016-14
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBYD6

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: BP013310.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.328	21	24	35	rVB	115408	169177	1.99%	0.149%
2	3.775	95	100	110	rBV2	91574	156632	1.84%	0.138%
3	3.864	111	115	121	rVB	116119	140767	1.66%	0.124%
4	3.987	131	136	141	rVB	123993	187336	2.20%	0.165%
5	4.087	149	153	157	rVB	49229	70923	0.83%	0.063%
6	4.246	178	180	187	rVB6	19233	29163	0.34%	0.026%
7	4.464	214	217	221	rVB3	27419	36274	0.43%	0.032%
8	4.605	237	241	245	rVB4	13933	18587	0.22%	0.016%
9	4.652	245	249	253	rVV	47831	67263	0.79%	0.059%
10	4.699	253	257	268	rVB	159582	233392	2.75%	0.206%
11	5.017	305	311	320	rBV3	388421	701712	8.26%	0.619%
12	5.122	327	329	333	rVB4	17247	20697	0.24%	0.018%
13	5.622	411	414	418	rVB	15952	20452	0.24%	0.018%
14	6.540	565	570	575	rVV	47177	68909	0.81%	0.061%
15	6.611	575	582	590	rVB	469498	708576	8.34%	0.625%
16	6.922	630	635	638	rBV5	18661	29634	0.35%	0.026%
17	7.099	655	665	679	rBV	1464948	2463525	28.99%	2.174%
18	7.263	687	693	703	rVB	1343194	2095137	24.65%	1.849%
19	7.369	707	711	717	rVV	54685	86451	1.02%	0.076%
20	7.469	719	728	737	rBV	1629594	2569173	30.23%	2.268%
21	7.546	737	741	750	rVB2	31559	56058	0.66%	0.049%
22	7.752	772	776	785	rVB4	14810	29577	0.35%	0.026%
23	7.940	801	808	815	rBV	2529503	3848176	45.28%	3.397%
24	8.152	840	844	850	rBV9	13818	30890	0.36%	0.027%
25	8.652	920	929	939	rBV	1163024	1854107	21.82%	1.637%
26	8.769	943	949	959	rVB	119876	202518	2.38%	0.179%
27	9.110	999	1007	1014	rBV	1488987	2489055	29.29%	2.197%
28	9.263	1029	1033	1038	rVB7	13116	22597	0.27%	0.020%
29	9.505	1068	1074	1086	rVB3	62523	119997	1.41%	0.106%
30	9.834	1119	1130	1137	rBV	1227939	2033151	23.92%	1.795%
31	9.987	1150	1156	1163	rBV6	17145	45124	0.53%	0.040%
32	10.052	1163	1167	1178	rVB6	28557	63844	0.75%	0.056%
33	10.187	1187	1190	1197	rVB5	11537	21678	0.26%	0.019%
34	10.369	1214	1221	1233	rBV	1729957	2958793	34.81%	2.612%
35	10.752	1277	1286	1293	rBV	3176289	5271313	62.02%	4.653%
36	10.893	1299	1310	1329	rVB	1100133	2031185	23.90%	1.793%

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Integrator: RTE

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Stop Thrs : 0

Peak Location: TOP

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

37	11.281	1372	1376	1384	rVV	11647	26550	0.31%	0.023%
38	11.540	1416	1420	1431	rVB8	14696	35831	0.42%	0.032%
39	12.004	1495	1499	1506	rVB9	11035	18549	0.22%	0.016%
40	12.334	1550	1555	1564	rVB2	35547	62581	0.74%	0.055%
41	12.857	1636	1644	1650	rBV8	16748	34244	0.40%	0.030%
42	13.104	1681	1686	1692	rVB3	29923	44210	0.52%	0.039%
43	13.204	1697	1703	1708	rBV9	11934	24468	0.29%	0.022%
44	13.393	1732	1735	1741	rVB2	31438	46692	0.55%	0.041%
45	13.469	1743	1748	1752	rBV8	17528	31738	0.37%	0.028%
46	13.763	1794	1798	1802	rVB4	12444	18595	0.22%	0.016%
47	13.857	1809	1814	1818	rBV5	25094	60690	0.71%	0.054%
48	13.904	1818	1822	1828	rVB2	117555	171283	2.02%	0.151%
49	13.987	1828	1836	1846	rBV	2845799	3979025	46.82%	3.512%
50	14.104	1851	1856	1862	rVB8	28005	49652	0.58%	0.044%
51	14.275	1872	1885	1901	rBV	3249522	4817647	56.69%	4.252%
52	14.463	1913	1917	1924	rVB3	43651	65704	0.77%	0.058%
53	14.540	1924	1930	1931	rBV6	13646	19726	0.23%	0.017%
54	14.581	1931	1937	1945	rVV	3708848	5693619	66.99%	5.026%
55	14.645	1945	1948	1955	rVB9	32466	66841	0.79%	0.059%
56	14.787	1965	1972	1986	rBV	599734	1180260	13.89%	1.042%
57	14.940	1994	1998	2002	rVB2	38032	54340	0.64%	0.048%
58	14.998	2002	2008	2013	rBV7	68996	126453	1.49%	0.112%
59	15.245	2047	2050	2056	rVB8	12440	19489	0.23%	0.017%
60	15.363	2066	2070	2078	rBV9	11706	30453	0.36%	0.027%
61	15.492	2086	2092	2099	rBV5	38663	55737	0.66%	0.049%
62	15.575	2099	2106	2113	rBV	3803968	5501722	64.74%	4.856%
63	15.628	2113	2115	2122	rVB4	25266	35774	0.42%	0.032%
64	15.698	2122	2127	2135	rBV	640334	857111	10.09%	0.757%
65	15.822	2142	2148	2153	rVB6	23757	38914	0.46%	0.034%
66	15.945	2162	2169	2176	rBV	287982	402634	4.74%	0.355%
67	16.039	2179	2185	2192	rBV	19915	52442	0.62%	0.046%
68	16.145	2199	2203	2210	rVB7	37242	57946	0.68%	0.051%
69	16.210	2210	2214	2219	rBV4	35931	48485	0.57%	0.043%
70	16.287	2223	2227	2234	rVB9	18026	33200	0.39%	0.029%
71	16.722	2298	2301	2306	rVB5	23887	32951	0.39%	0.029%
72	16.875	2323	2327	2334	rVB	54090	75859	0.89%	0.067%
73	17.081	2357	2362	2366	rVB8	14827	25758	0.30%	0.023%
74	17.222	2381	2386	2389	rBV7	26944	44282	0.52%	0.039%
75	17.345	2400	2407	2417	rBV	3980760	5877448	69.16%	5.188%
76	17.439	2417	2423	2432	rVV	3504868	5151818	60.62%	4.547%

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77	17.534	2436	2439	2446	rVB7	39837	65682	0.77%	0.058%
78	17.722	2469	2471	2481	rVB7	18338	36300	0.43%	0.032%
79	17.998	2515	2518	2524	rVB7	34357	45973	0.54%	0.041%
80	18.098	2530	2535	2541	rBV6	35941	68937	0.81%	0.061%
81	18.210	2550	2554	2564	rBV	421483	610702	7.19%	0.539%
82	18.545	2607	2611	2617	rBV3	75562	124830	1.47%	0.110%
83	18.628	2622	2625	2628	rVB3	33652	35605	0.42%	0.031%
84	18.828	2656	2659	2660	rBV3	28075	30374	0.36%	0.027%
85	18.969	2678	2683	2691	rBV5	183106	291675	3.43%	0.257%
86	19.251	2728	2731	2737	rVB6	64668	98335	1.16%	0.087%
87	19.328	2738	2744	2746	rBV	75837	150235	1.77%	0.133%
88	19.445	2760	2764	2768	rBV2	132702	197804	2.33%	0.175%
89	19.675	2798	2803	2816	rVB	5423949	7129434	83.89%	6.293%
90	20.180	2887	2889	2895	rVB2	94680	133471	1.57%	0.118%
91	20.616	2960	2963	2967	rVB	188807	235462	2.77%	0.208%
92	20.669	2969	2972	2976	rVB2	118890	150578	1.77%	0.133%
93	21.127	3046	3050	3059	rVB	896795	1254131	14.76%	1.107%
94	21.433	3097	3102	3107	rBV	5647251	7141480	84.03%	6.304%
95	22.074	3208	3211	3223	rVB3	363692	610457	7.18%	0.539%
96	23.127	3384	3390	3396	rBV	2402049	3973613	46.75%	3.507%
97	23.751	3490	3496	3509	rVB2	3903863	8498821	100.00%	7.502%
98	23.916	3517	3524	3532	rVB2	3700868	7696517	90.56%	6.793%
99	24.527	3622	3628	3637	rVB	2748499	5698682	67.05%	5.030%
100	24.633	3640	3646	3663	rVB2	1055085	3090856	36.37%	2.728%

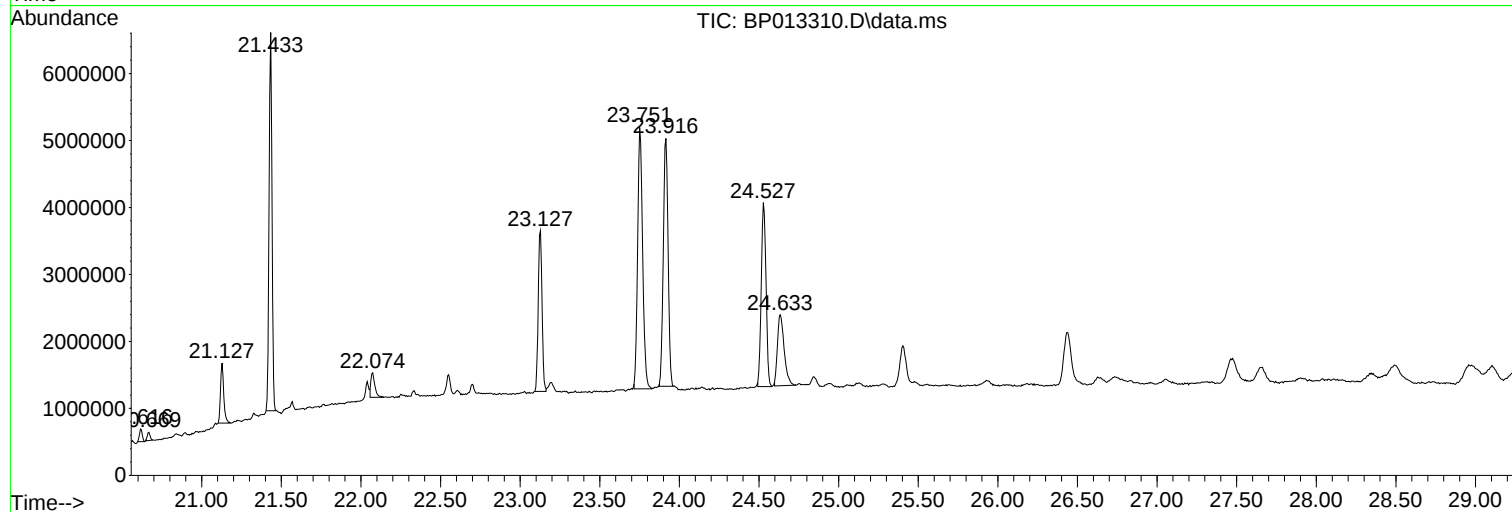
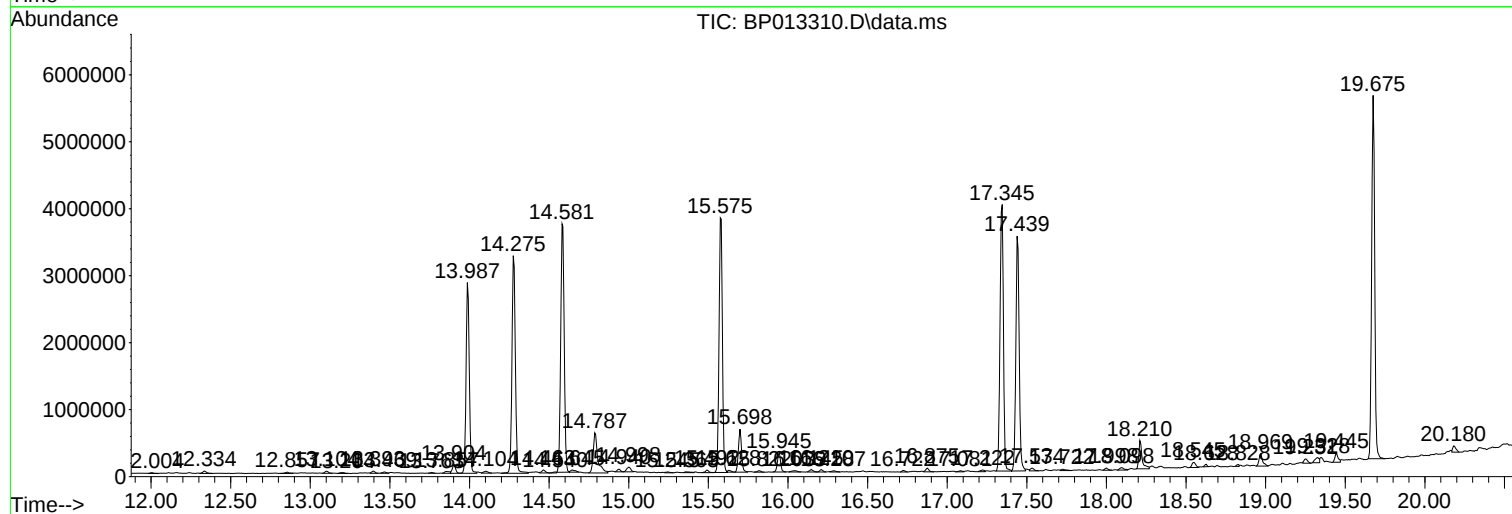
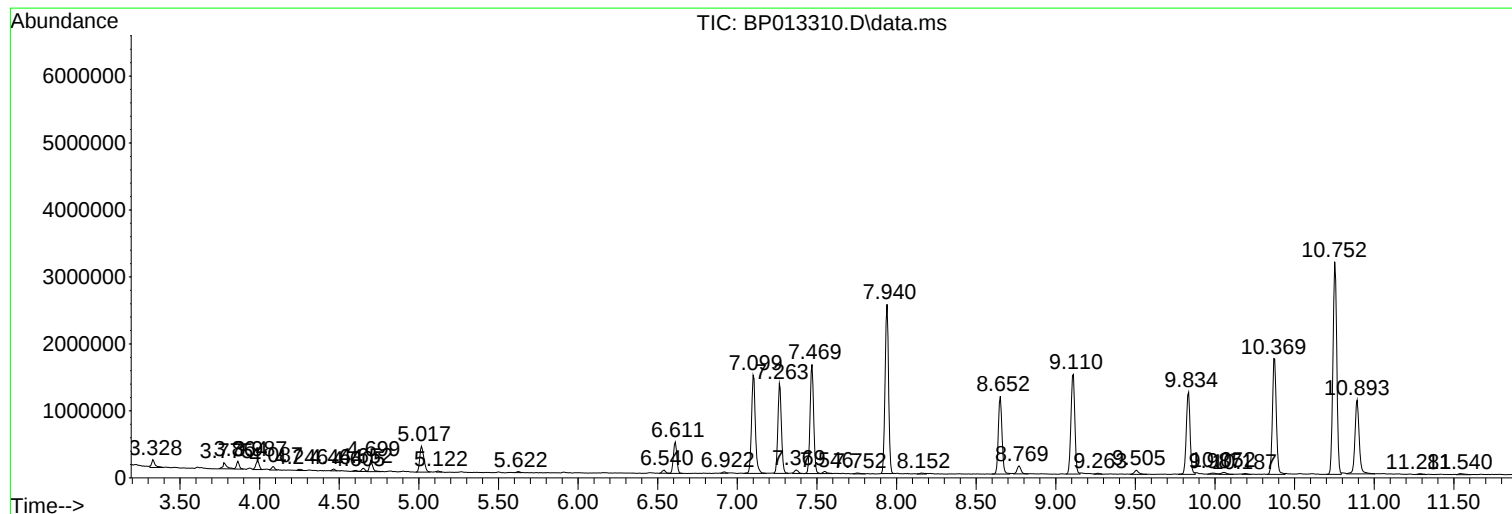
Sum of corrected areas: 113292518

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Instrument :
BNA_P
ClientSampleId :
DBYD6

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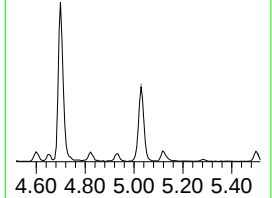
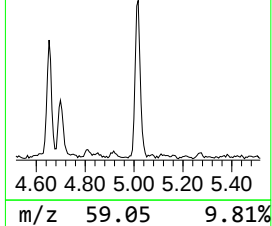
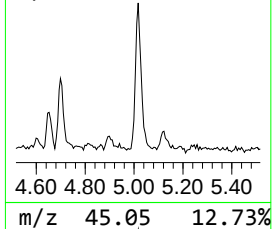
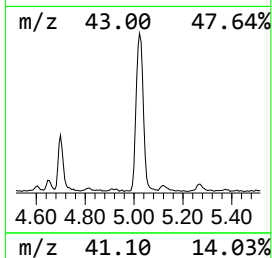
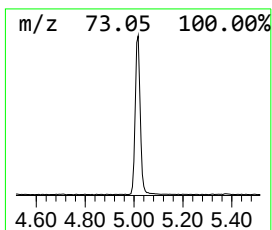
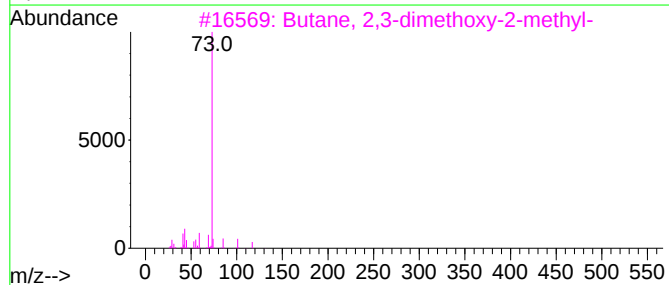
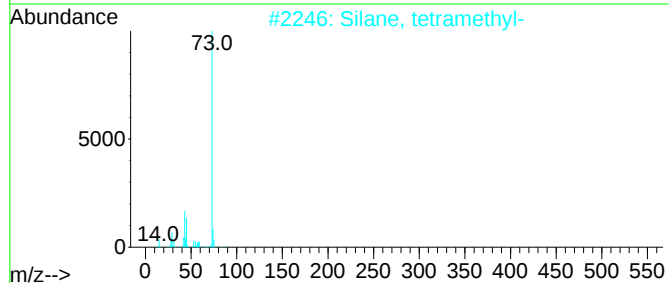
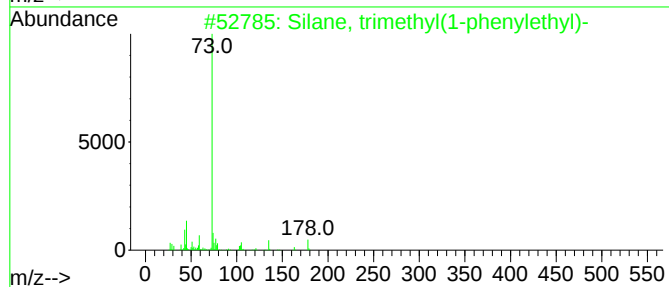
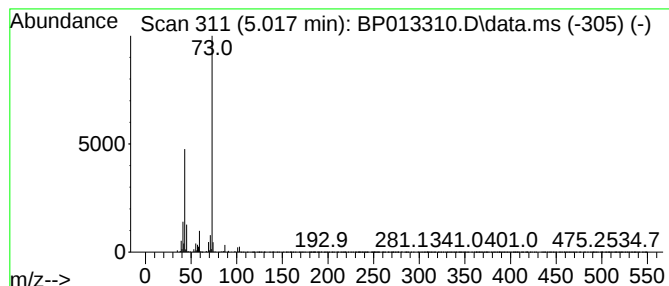
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.017	3.65 ng/ul	701712	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silane, trimethyl(1-phenylethyl)-	178	C11H18Si	017961-78-3	38
2			Silane, tetramethyl-	88	C4H12Si	000075-76-3	37
3			Butane, 2,3-dimethoxy-2-methyl-	132	C7H16O2	074421-00-4	33
4			Silane, 2-butenyltrimethyl-, (E)-	128	C7H16Si	017486-12-3	33
5			1,3-Dioxolane, 2-(1-methylethyl)-	116	C6H12O2	000822-83-3	25



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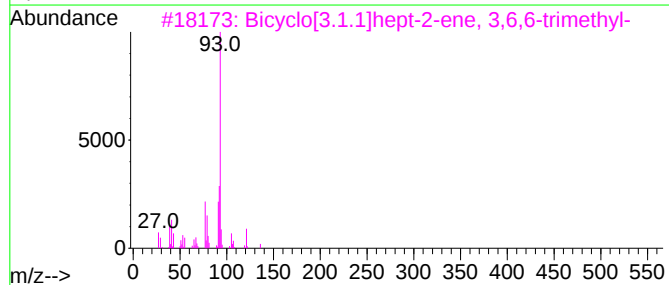
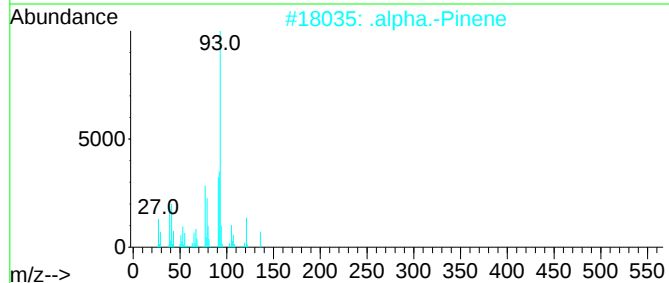
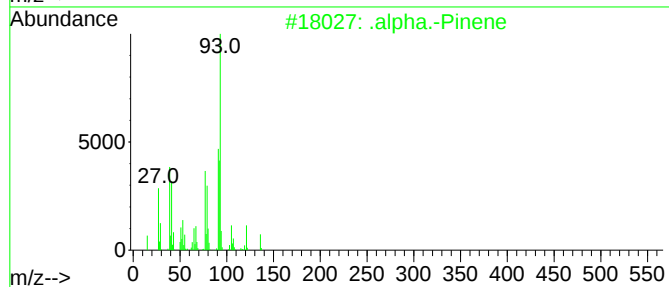
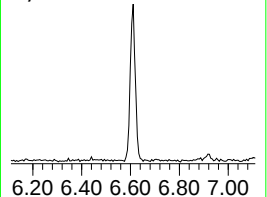
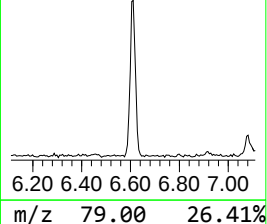
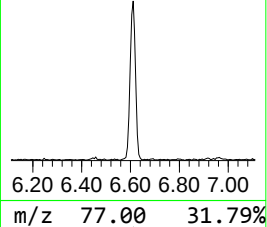
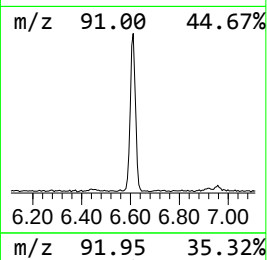
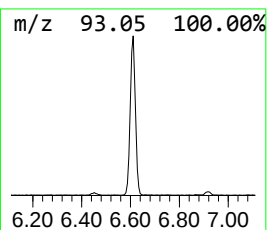
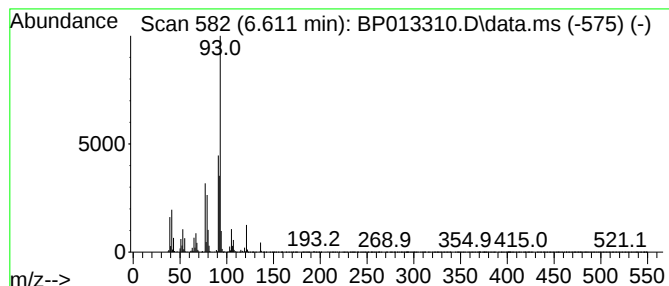
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 .alpha.-Pinene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.611	3.68 ng/ul	708576	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.		.alpha.-Pinene	136	C10H16	000080-56-8	95
2	.		.alpha.-Pinene	136	C10H16	000080-56-8	94
3	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...			136	C10H16	004889-83-2	94
4	.gamma.-Terpinene			136	C10H16	000099-85-4	91
5	(1R)-2,6,6-Trimethylbicyclo[3.1....			136	C10H16	007785-70-8	91



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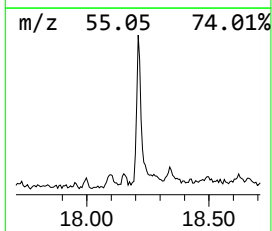
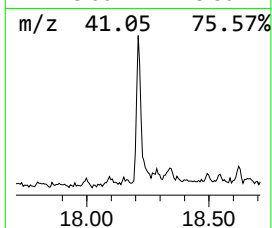
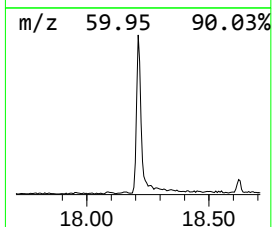
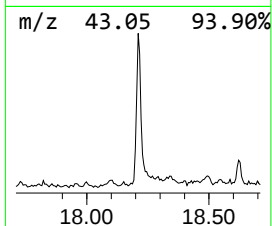
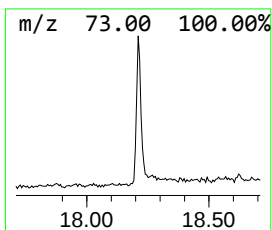
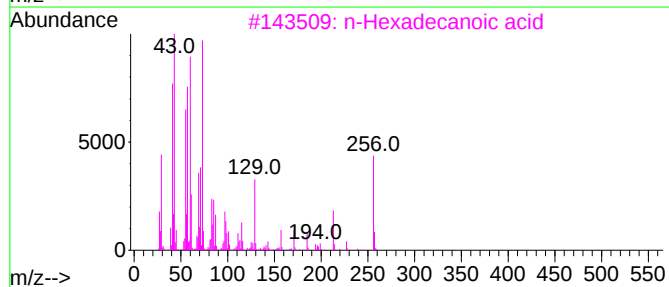
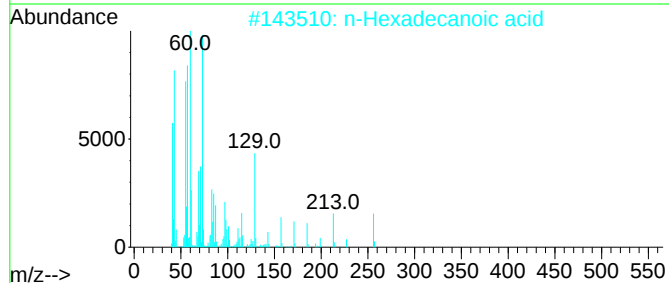
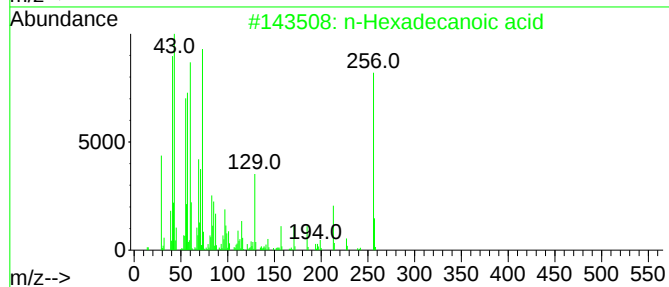
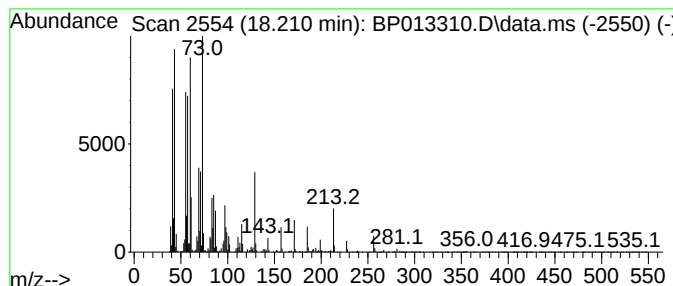
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.210	2.08 ng/ul	610702	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	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	236	C16H32O2	000057-10-3	97
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013310.D
Acq On : 24 Dec 2022 21:28
Operator : CG/JU
Sample : N6016-14
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD6

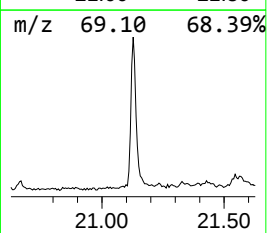
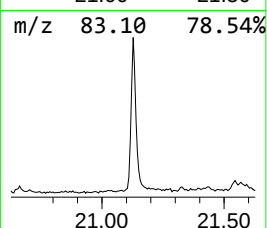
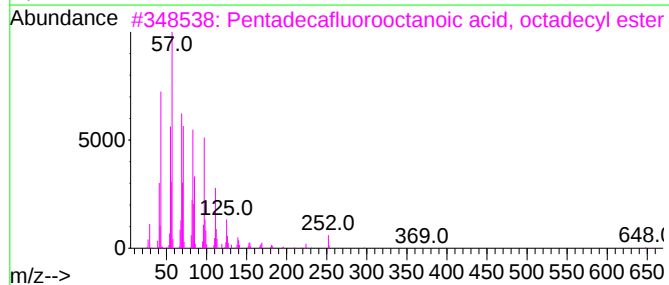
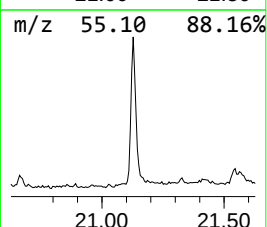
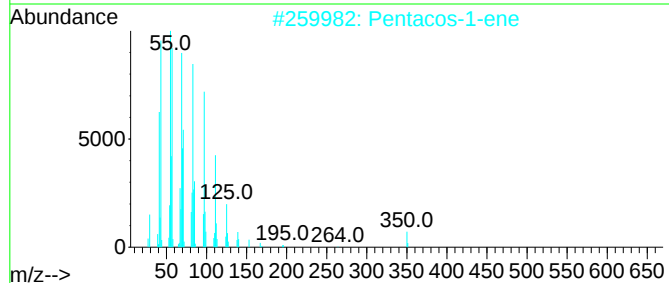
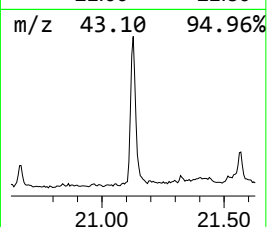
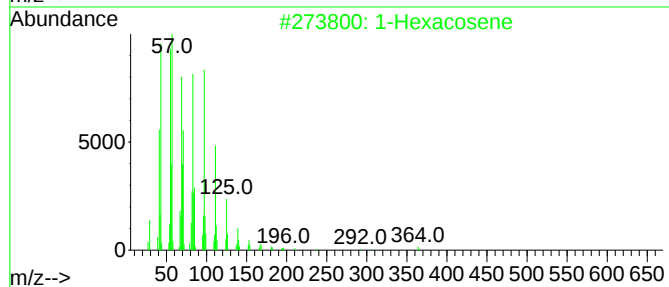
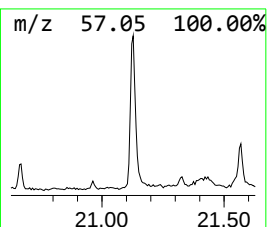
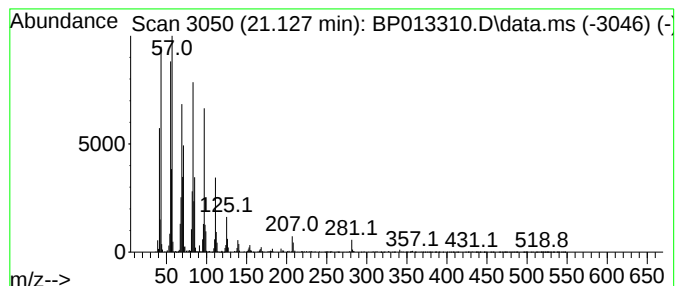
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.128	3.51 ng/ul	1254130	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene			364	C26H52	018835-33-1	94
2	Pentacos-1-ene			350	C25H50	016980-85-1	94
3	Pentadecafluorooctanoic acid, oc...			666	C26H37F15O2	1000400-04-8	93
4	Pentafluoropropionic acid, octad...			416	C21H37F5O2	959261-25-7	91
5	Heptafluorobutyric acid, n-octad...			466	C22H37F7O2	000400-57-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013310.D
Acq On : 24 Dec 2022 21:28
Operator : CG/JU
Sample : N6016-14
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD6

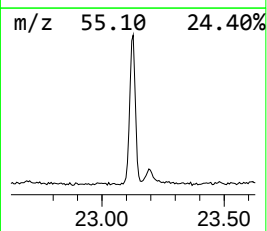
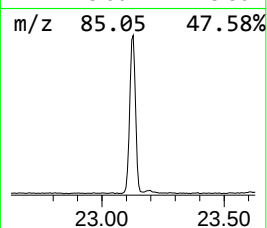
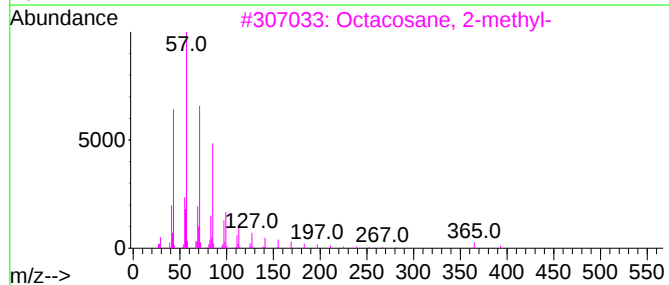
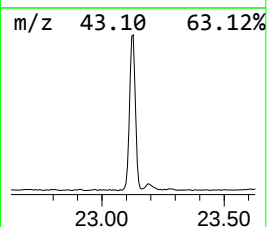
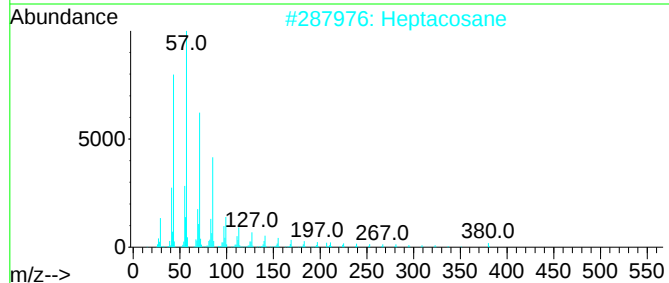
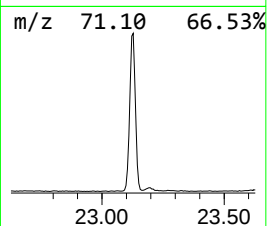
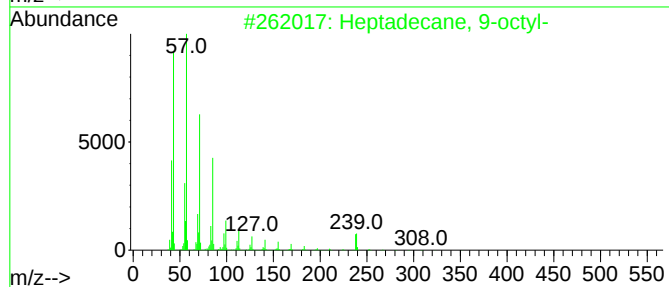
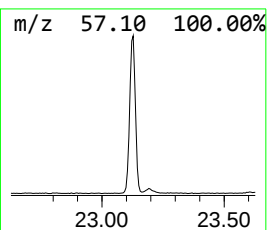
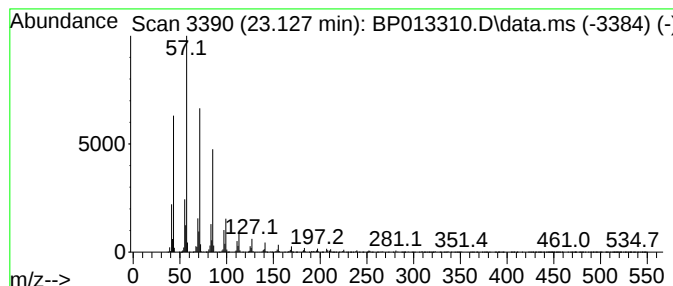
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.127	10.33 ng/ul	3973610	Perylene-d12	23.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2			Heptacosane	380	C27H56	000593-49-7	91
3			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4			Hexacosane	366	C26H54	000630-01-3	91
5			Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013310.D
Acq On : 24 Dec 2022 21:28
Operator : CG/JU
Sample : N6016-14
Misc :
ALS Vial : 22 Sample Multiplier: 1

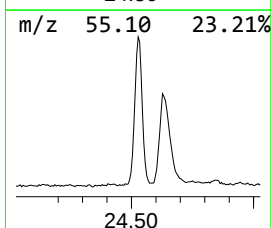
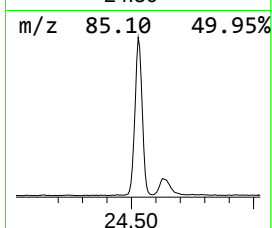
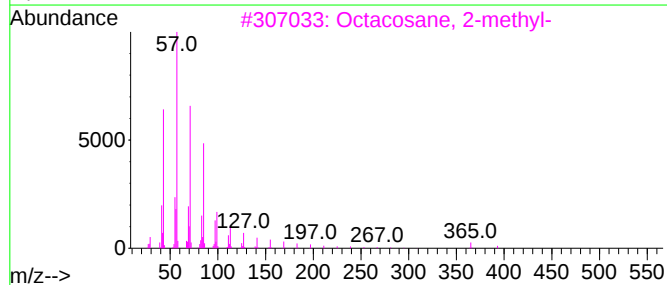
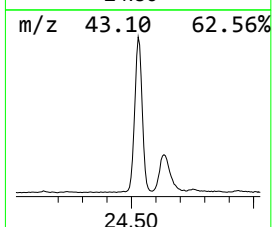
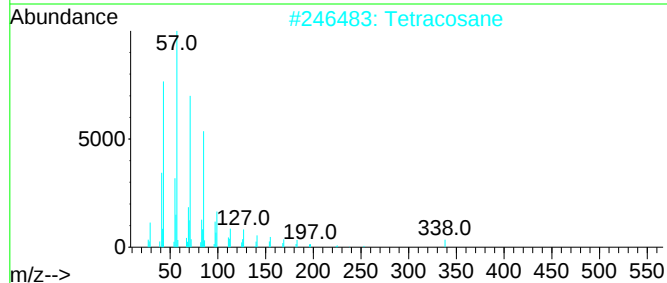
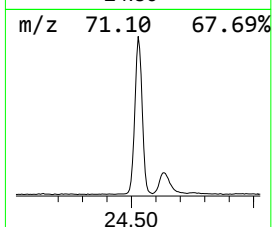
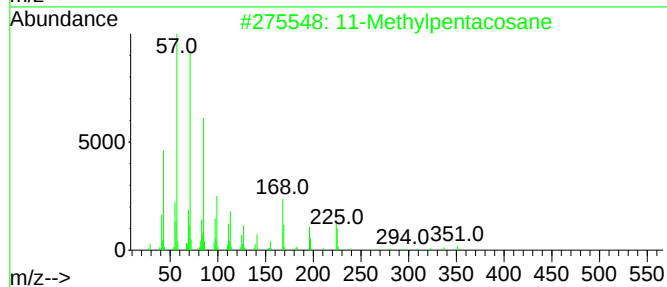
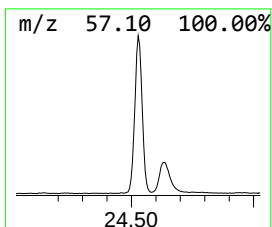
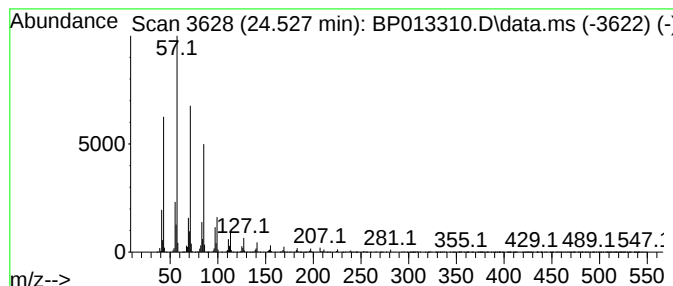
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
24.527	14.81 ng/ul	5698680	Perylene-d12		23.916	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	11-Methylpentacosane		366	C26H54	015689-71-1	94
2	Tetracosane		338	C24H50	000646-31-1	93
3	Octacosane, 2-methyl-		408	C29H60	001560-98-1	91
4	Heneicosane		296	C21H44	000629-94-7	91
5	Hexatriacontane		507	C36H74	000630-06-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013310.D
Acq On : 24 Dec 2022 21:28
Operator : CG/JU
Sample : N6016-14
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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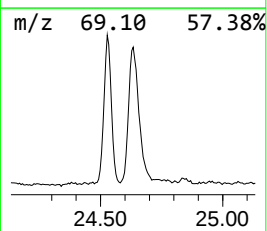
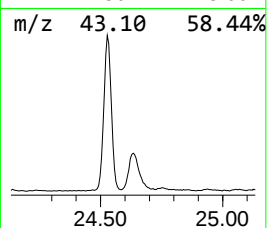
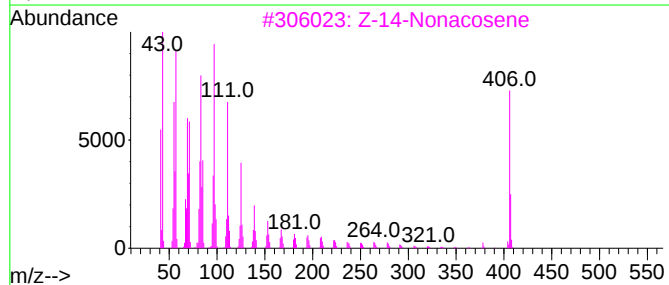
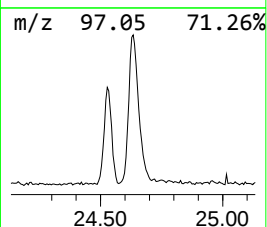
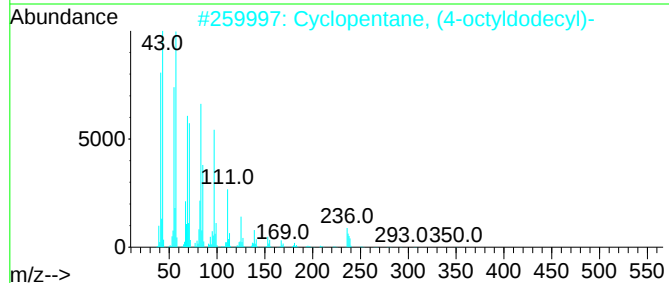
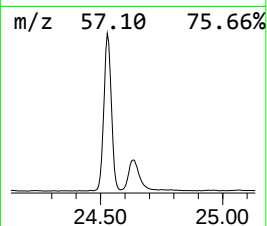
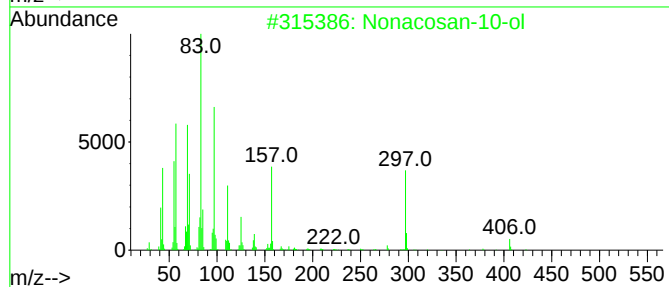
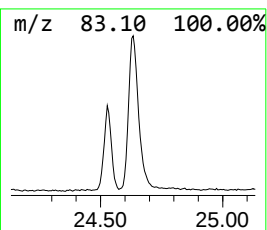
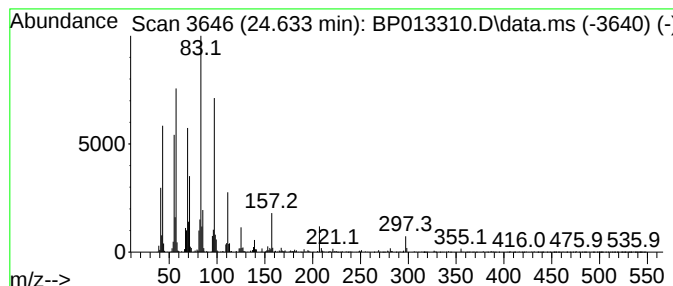
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Nonacosan-10-ol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.633	8.03 ng/ul	3090860	Perylene-d12	23.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonacosan-10-ol	424	C29H60O	000504-55-2	86
2			Cyclopentane, (4-octyldodecyl)-	350	C25H50	005638-09-5	52
3			Z-14-Nonacosene	406	C29H58	1010131-18-9	50
4			Hexacosyl heptafluorobutyrate	578	C30H53F7O2	1000351-83-3	50
5			Hexacosyl pentafluoropropionate	528	C29H53F5O2	1000351-81-2	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013310.D
Acq On : 24 Dec 2022 21:28
Operator : CG/JU
Sample : N6016-14
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD6

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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.alpha.-Pinene	6.611	3.7	ng/ul	708576	1	7.940	3848180	20.0
n-Hexadecanoic ...	18.210	2.1	ng/ul	610702	4	17.345	5877450	20.0
(DEL) Alkane: S...	21.128	3.5	ng/ul	1254130	5	21.433	7141480	20.0
(DEL) Alkane: S...	23.127	10.3	ng/ul	3973610	6	23.916	7696520	20.0
(DEL) Alkane: S...	24.527	14.8	ng/ul	5698680	6	23.916	7696520	20.0
Nonacosan-10-ol	24.633	8.0	ng/ul	3090860	6	23.916	7696520	20.0