

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
 Data File : BP013325.D
 Acq On : 25 Dec 2022 05:54
 Operator : CG/JU
 Sample : N6018-18
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBYK3

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: BP013325.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	31	rVB	123049	153327	1.46%	0.109%
2	3.758	94	97	113	rBV	915564	1361614	13.00%	0.966%
3	4.081	148	152	156	rBV	41499	57474	0.55%	0.041%
4	5.028	309	313	319	rVB2	51969	77869	0.74%	0.055%
5	5.258	348	352	359	rVB	30668	47737	0.46%	0.034%
6	5.975	470	474	484	rVB9	12953	28414	0.27%	0.020%
7	6.452	548	555	560	rVB4	31895	60309	0.58%	0.043%
8	6.611	575	582	591	rVB	3667457	5624369	53.69%	3.989%
9	6.917	629	634	640	rBV3	63361	113832	1.09%	0.081%
10	7.099	657	665	676	rBV	1989805	3201928	30.56%	2.271%
11	7.264	687	693	704	rVB	1663733	2509970	23.96%	1.780%
12	7.369	705	711	721	rVV	771554	1233198	11.77%	0.875%
13	7.469	721	728	735	rVV	2095195	3253789	31.06%	2.308%
14	7.940	800	808	815	rBV	2234891	3519958	33.60%	2.496%
15	8.005	815	819	824	rVB4	15329	24596	0.23%	0.017%
16	8.158	838	845	850	rBV	50340	79820	0.76%	0.057%
17	8.211	850	854	862	rVB3	20295	38304	0.37%	0.027%
18	8.652	922	929	939	rBV	2267497	3640550	34.75%	2.582%
19	9.105	998	1006	1014	rBV	1914189	3135517	29.93%	2.224%
20	9.263	1028	1033	1039	rVB4	36012	56720	0.54%	0.040%
21	9.505	1067	1074	1084	rBV2	75772	153055	1.46%	0.109%
22	9.828	1121	1129	1141	rBV	1573566	2608526	24.90%	1.850%
23	10.052	1163	1167	1178	rVB5	56598	131064	1.25%	0.093%
24	10.369	1213	1221	1238	rBV	2390367	4050940	38.67%	2.873%
25	10.752	1277	1286	1293	rBV	2937321	4814605	45.96%	3.415%
26	10.893	1299	1310	1322	rVV	1341172	2366655	22.59%	1.678%
27	11.222	1363	1366	1371	rVB7	12233	21588	0.21%	0.015%
28	11.534	1412	1419	1426	rBV7	12947	31166	0.30%	0.022%
29	12.005	1492	1499	1503	rVB8	9624	22217	0.21%	0.016%
30	12.157	1521	1525	1531	rVB	45637	58635	0.56%	0.042%
31	12.334	1551	1555	1559	rVV2	36244	51747	0.49%	0.037%
32	12.781	1626	1631	1636	rVV9	14121	33329	0.32%	0.024%
33	13.110	1682	1687	1691	rBV4	14430	23671	0.23%	0.017%
34	13.340	1722	1726	1729	rVV6	15542	26182	0.25%	0.019%
35	13.393	1729	1735	1740	rVB	88119	128257	1.22%	0.091%
36	13.451	1740	1745	1750	rBV7	29218	51415	0.49%	0.036%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
 Data File : BP013325.D
 Acq On : 25 Dec 2022 05:54
 Operator : CG/JU
 Sample : N6018-18
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBYK3

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

37	13.504	1750	1754	1757	rBV2	29448	39141	0.37%	0.028%
38	13.899	1817	1821	1827	rBV4	88005	131325	1.25%	0.093%
39	13.987	1828	1836	1844	rVV	3665663	4994296	47.67%	3.542%
40	14.051	1844	1847	1851	rVV4	40575	62501	0.60%	0.044%
41	14.099	1852	1855	1864	rVB10	32707	62049	0.59%	0.044%
42	14.275	1878	1885	1898	rBV2	4248461	6221966	59.39%	4.413%
43	14.416	1903	1909	1912	rVV6	36910	69467	0.66%	0.049%
44	14.463	1914	1917	1923	rVV	41656	64943	0.62%	0.046%
45	14.540	1926	1930	1931	rBV4	20814	27318	0.26%	0.019%
46	14.581	1931	1937	1944	rVV	3765517	5455667	52.08%	3.869%
47	14.646	1944	1948	1961	rVB6	83340	245261	2.34%	0.174%
48	14.787	1966	1972	1976	rBV	919786	1449659	13.84%	1.028%
49	14.822	1976	1978	1988	rVB3	218181	347835	3.32%	0.247%
50	14.934	1993	1997	2002	rVV2	68026	113498	1.08%	0.080%
51	14.987	2002	2006	2011	rVV5	56711	108949	1.04%	0.077%
52	15.034	2011	2014	2019	rVV5	41140	70768	0.68%	0.050%
53	15.110	2023	2027	2030	rVV6	17586	30335	0.29%	0.022%
54	15.304	2056	2060	2065	rVB5	17612	22642	0.22%	0.016%
55	15.369	2066	2071	2074	rBV5	15196	27898	0.27%	0.020%
56	15.416	2077	2079	2084	rVB5	17362	22871	0.22%	0.016%
57	15.493	2084	2092	2099	rBV2	149390	238827	2.28%	0.169%
58	15.575	2099	2106	2113	rBV	5005725	7254538	69.25%	5.145%
59	15.698	2121	2127	2134	rBV	849421	1157788	11.05%	0.821%
60	15.822	2143	2148	2154	rVB5	39842	68625	0.66%	0.049%
61	15.940	2165	2168	2176	rVB	92536	128874	1.23%	0.091%
62	16.022	2178	2182	2184	rBV2	58424	85049	0.81%	0.060%
63	16.151	2198	2204	2210	rVB2	118889	160251	1.53%	0.114%
64	16.210	2210	2214	2217	rBV5	16678	29053	0.28%	0.021%
65	16.293	2224	2228	2233	rVB7	18921	34293	0.33%	0.024%
66	16.469	2256	2258	2262	rVB	20611	24707	0.24%	0.018%
67	16.728	2297	2302	2307	rBV4	64668	102637	0.98%	0.073%
68	16.840	2317	2321	2326	rVB3	49997	68555	0.65%	0.049%
69	17.087	2359	2363	2366	rBV6	16888	26652	0.25%	0.019%
70	17.128	2366	2370	2374	rVV5	28041	39743	0.38%	0.028%
71	17.222	2383	2386	2389	rBV5	17522	23971	0.23%	0.017%
72	17.257	2389	2392	2395	rVV5	19248	24774	0.24%	0.018%
73	17.340	2400	2406	2417	rBV	3912945	5693212	54.35%	4.038%
74	17.440	2417	2423	2431	rVV	5094624	6891476	65.78%	4.888%
75	17.551	2438	2442	2445	rBV6	23383	33518	0.32%	0.024%
76	17.616	2450	2453	2459	rVB6	55409	78213	0.75%	0.055%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
 Data File : BP013325.D
 Acq On : 25 Dec 2022 05:54
 Operator : CG/JU
 Sample : N6018-18
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBYK3

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

77	17.692	2463	2466	2469	rBV4	40444	58295	0.56%	0.041%
78	17.998	2515	2518	2522	rVB6	34447	41157	0.39%	0.029%
79	18.098	2530	2535	2539	rBV7	39374	76562	0.73%	0.054%
80	18.157	2541	2545	2550	rBV3	67644	103787	0.99%	0.074%
81	18.210	2550	2554	2565	rBV	1056503	1602516	15.30%	1.137%
82	18.545	2608	2611	2616	rVB6	74351	99469	0.95%	0.071%
83	18.645	2625	2628	2634	rVB	82700	112267	1.07%	0.080%
84	19.104	2703	2706	2711	rBV3	163770	202583	1.93%	0.144%
85	19.345	2737	2747	2752	rBV3	247210	707370	6.75%	0.502%
86	19.445	2759	2764	2769	rBV2	323225	502672	4.80%	0.357%
87	19.675	2797	2803	2812	rBV	6836339	8948699	85.42%	6.347%
88	19.798	2822	2824	2826	rBV2	88079	100890	0.96%	0.072%
89	19.881	2835	2838	2843	rVB7	113643	141317	1.35%	0.100%
90	20.616	2959	2963	2967	rBV	427573	556102	5.31%	0.394%
91	21.122	3044	3049	3061	rBV2	2098132	4063708	38.79%	2.882%
92	21.433	3096	3102	3107	rBV	4995317	7173102	68.47%	5.087%
93	22.033	3202	3204	3207	rBV2	324420	404058	3.86%	0.287%
94	22.069	3207	3210	3220	rVB	626744	960400	9.17%	0.681%
95	23.122	3385	3389	3396	rBV	1161944	1801079	17.19%	1.277%
96	23.751	3490	3496	3510	rVB2	4768045	10475811	100.00%	7.430%
97	23.916	3517	3524	3531	rVB2	3374923	7020691	67.02%	4.979%
98	24.416	3603	3609	3618	rVB3	2088819	4265883	40.72%	3.025%
99	24.521	3623	3627	3636	rVB	1352318	2803598	26.76%	1.988%
100	24.627	3639	3645	3655	rBV2	1618175	4251109	40.58%	3.015%

Sum of corrected areas: 140998617

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013325.D
Acq On : 25 Dec 2022 05:54
Operator : CG/JU
Sample : N6018-18
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYK3

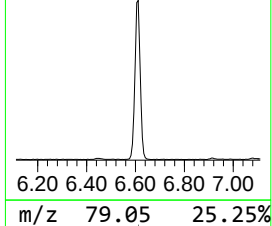
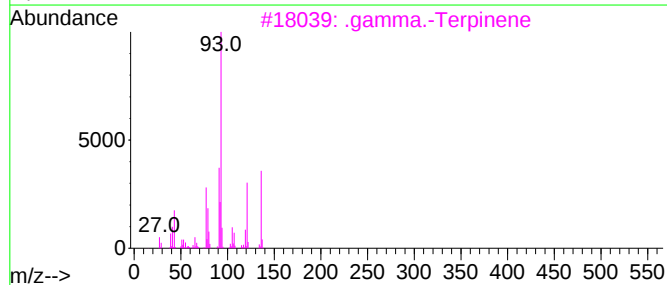
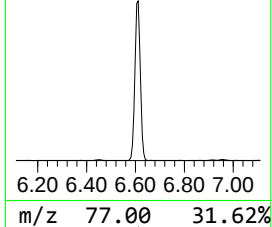
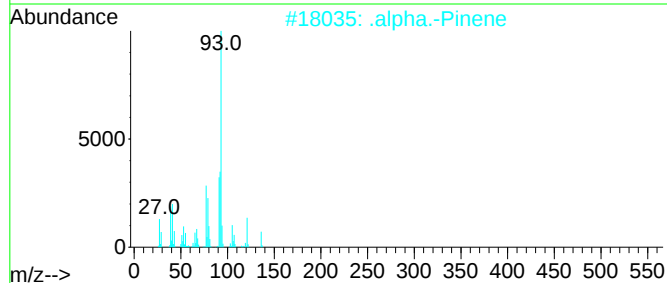
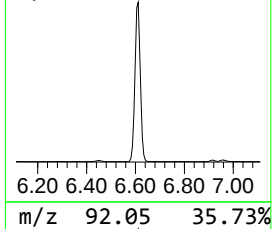
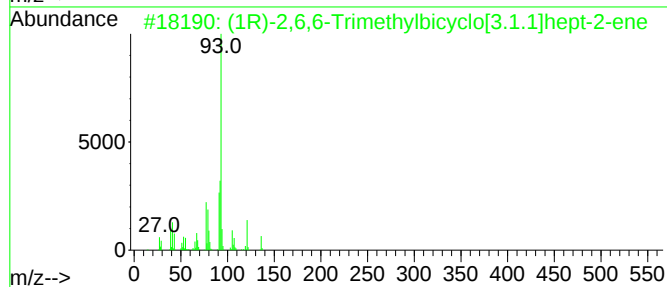
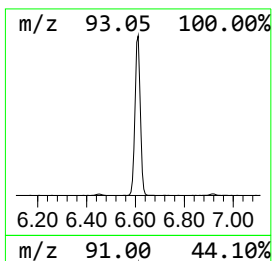
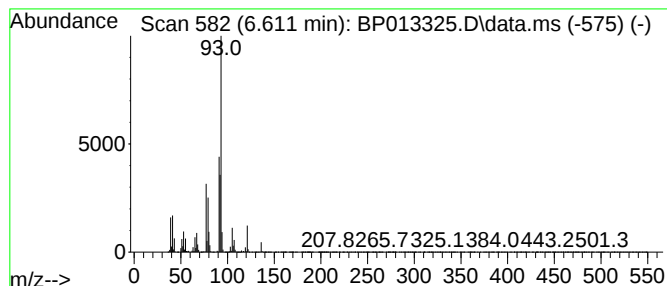
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.611	31.96 ng/ul	5624370	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-70-8	95		
2	.alpha.-Pinene	136	C10H16	000080-56-8	94		
3	.gamma.-Terpinene	136	C10H16	000099-85-4	91		
4	(1S)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-26-4	91		
5	.alpha.-Pinene	136	C10H16	000080-56-8	91		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013325.D
Acq On : 25 Dec 2022 05:54
Operator : CG/JU
Sample : N6018-18
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYK3

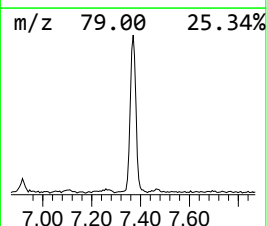
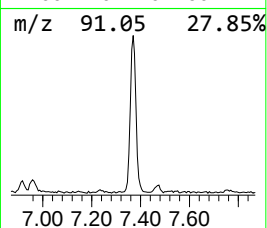
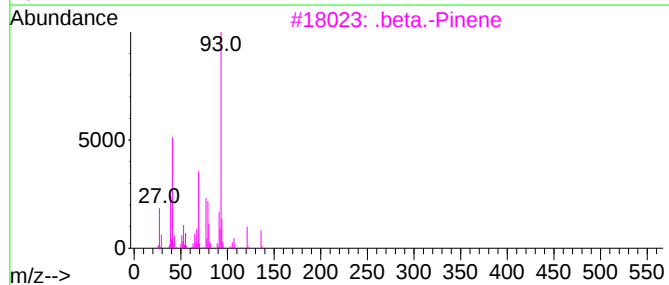
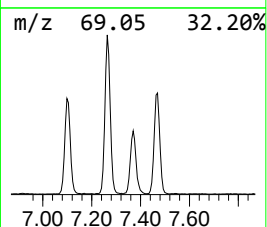
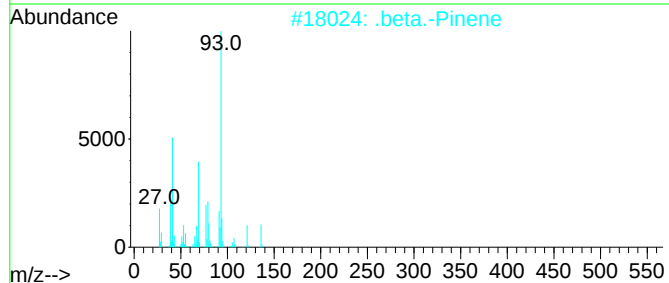
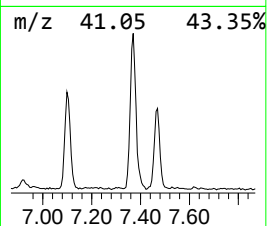
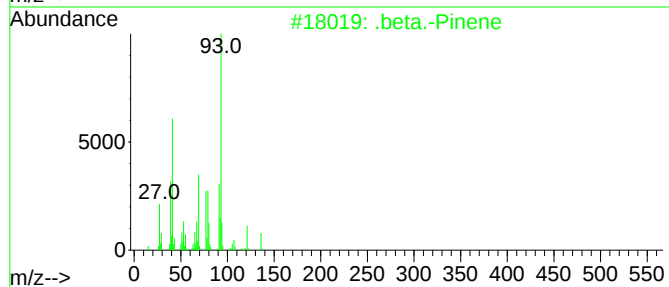
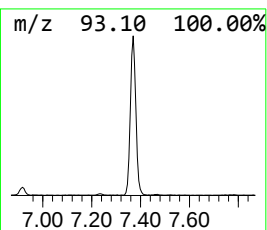
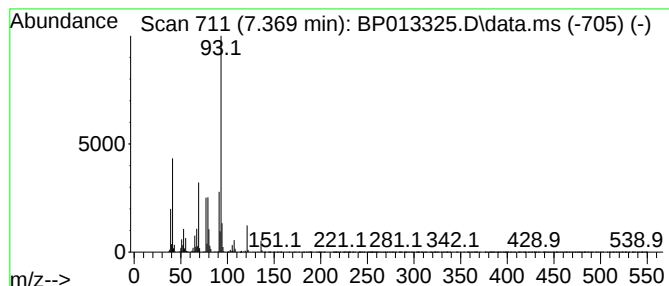
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 .beta.-Pinene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.369	7.01 ng/ul	1233200	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	beta.-Pinene	136	C10H16	000127-91-3	93	
2	.	beta.-Pinene	136	C10H16	000127-91-3	91	
3	.	beta.-Pinene	136	C10H16	000127-91-3	91	
4	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	91		
5	Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	91		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013325.D
Acq On : 25 Dec 2022 05:54
Operator : CG/JU
Sample : N6018-18
Misc :
ALS Vial : 37 Sample Multiplier: 1

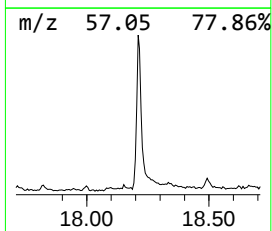
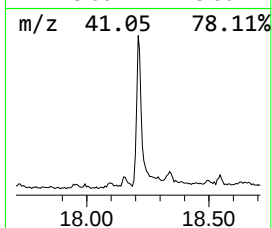
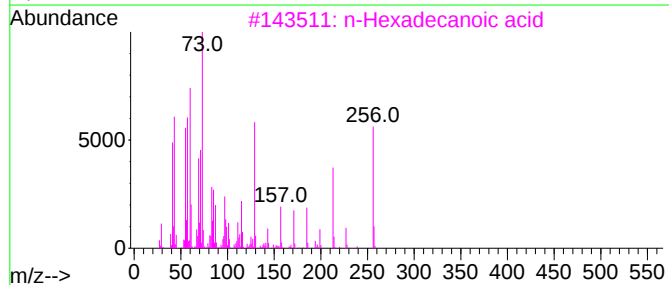
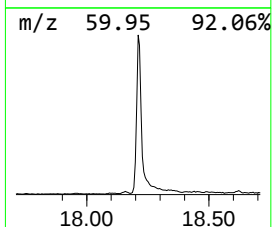
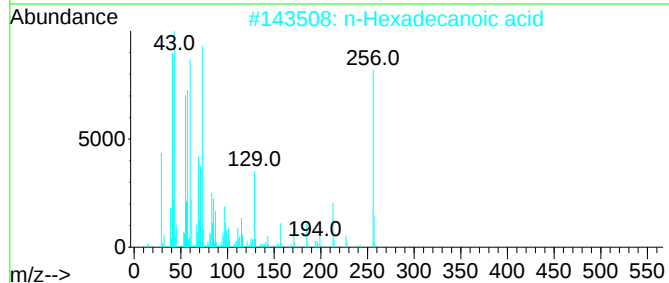
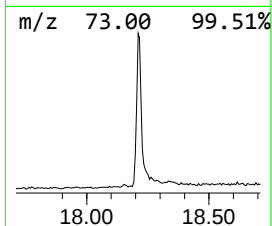
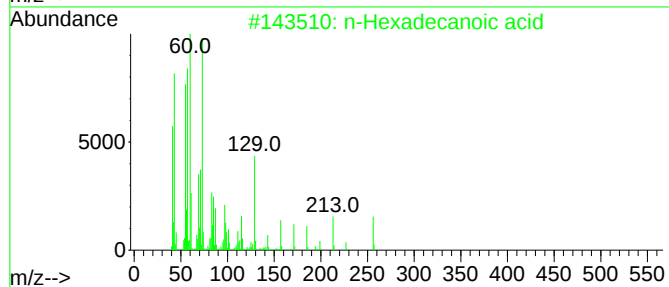
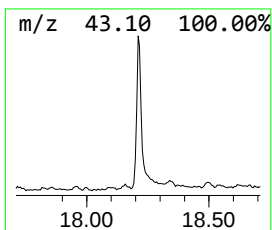
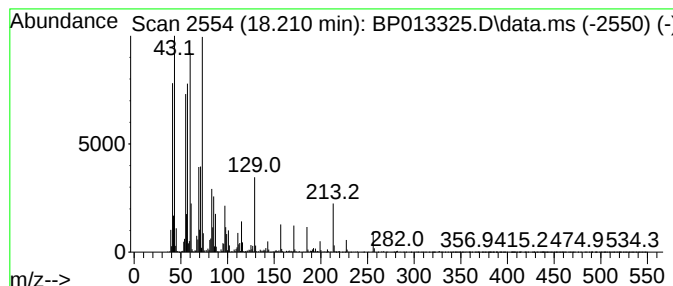
Instrument :
BNA_P
ClientSampleId :
DBYK3

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 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.210	5.63 ng/ul	1602520	Phenanthrene-d10		17.340	
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	97
4	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	96
5	Tridecanoic acid		214	C13H26O2	000638-53-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013325.D
Acq On : 25 Dec 2022 05:54
Operator : CG/JU
Sample : N6018-18
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYK3

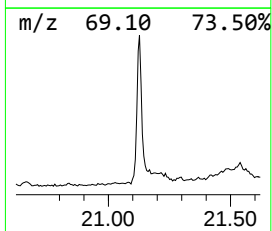
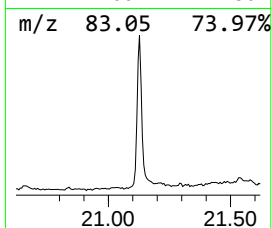
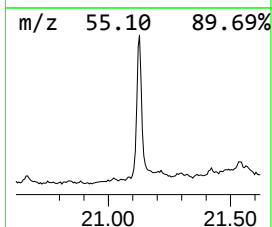
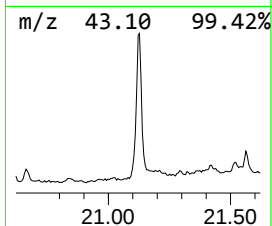
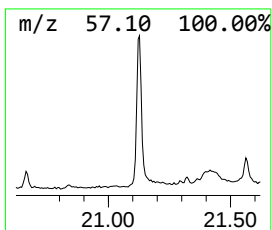
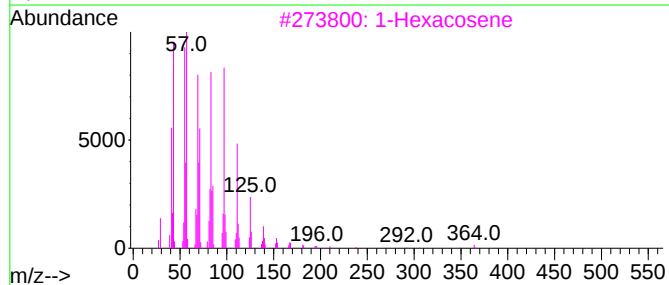
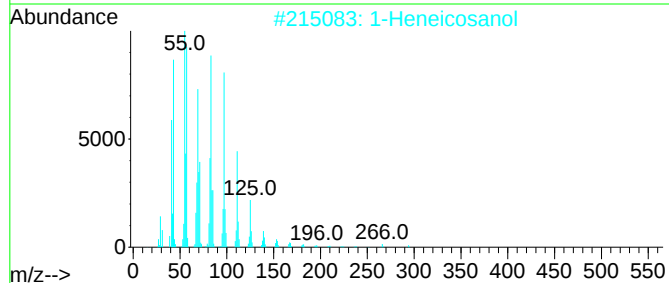
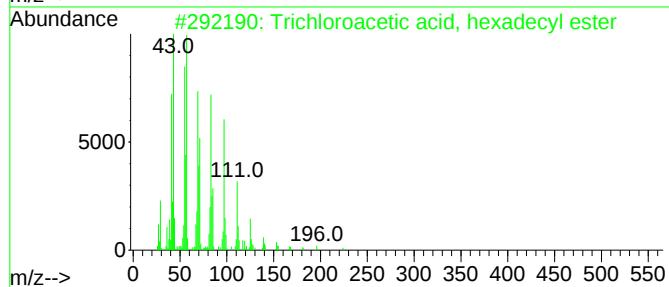
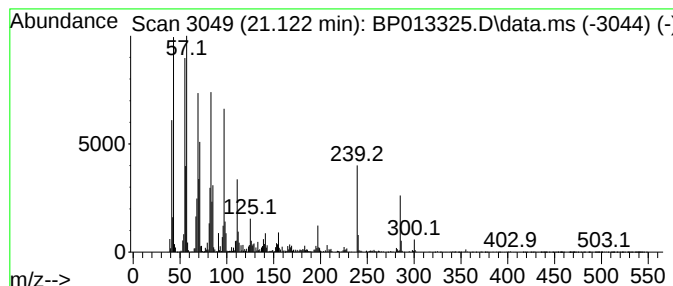
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Trichloroacetic acid, hexad... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.122	11.33 ng/ul	4063710	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	93
2			1-Heneicosanol	312	C21H44O	015594-90-8	93
3			1-Hexacosene	364	C26H52	018835-33-1	91
4			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
5			Heptacos-1-ene	378	C27H54	015306-27-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013325.D
Acq On : 25 Dec 2022 05:54
Operator : CG/JU
Sample : N6018-18
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYK3

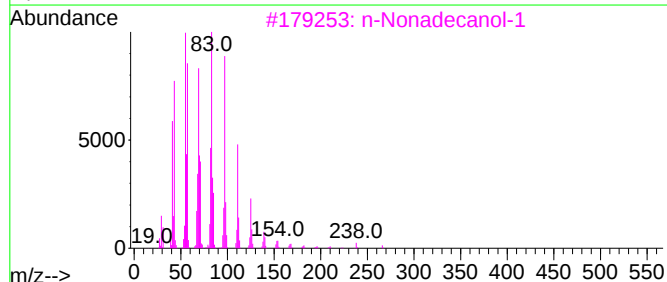
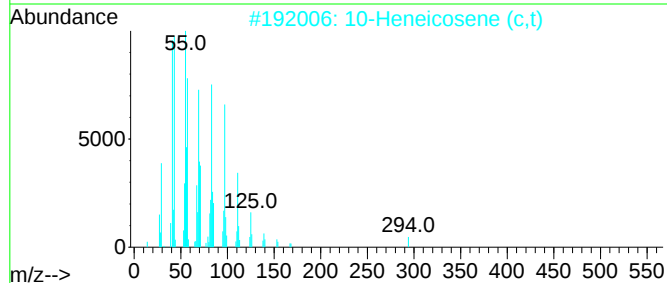
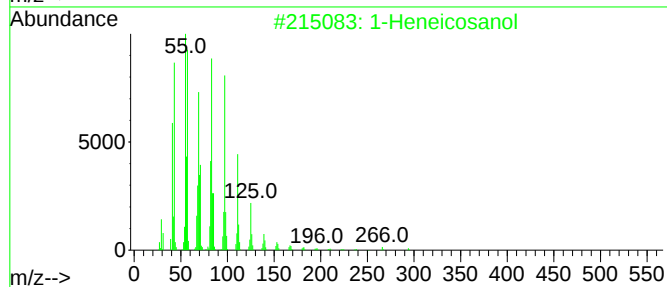
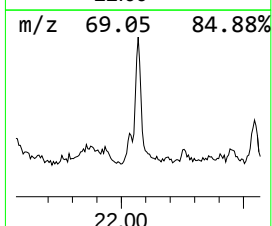
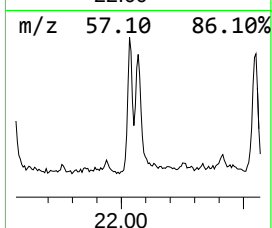
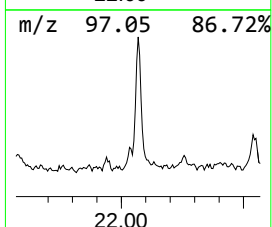
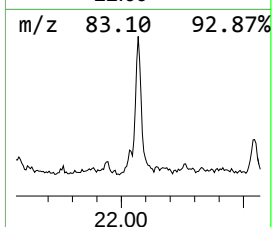
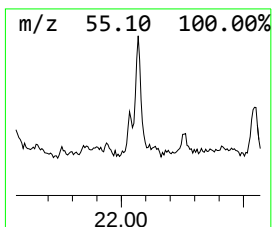
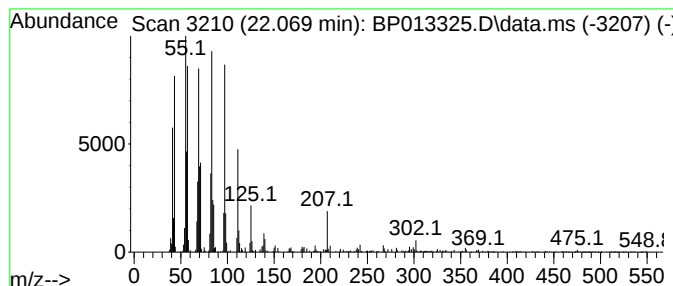
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 1-Heneicosanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.069	2.68 ng/ul	960400	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	95
2			10-Heneicosene (c,t)	294	C21H42	095008-11-0	95
3			n-Nonadecanol-1	284	C19H40O	001454-84-8	94
4			Behenic alcohol	326	C22H46O	000661-19-8	93
5			E-15-Heptadecenal	252	C17H32O	1000130-97-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013325.D
Acq On : 25 Dec 2022 05:54
Operator : CG/JU
Sample : N6018-18
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYK3

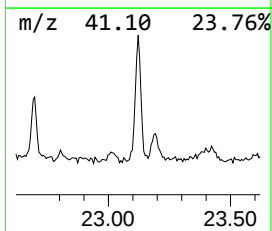
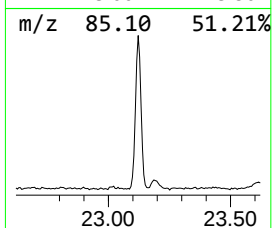
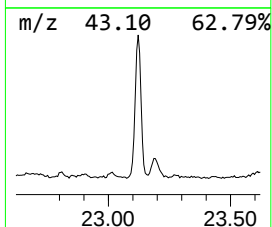
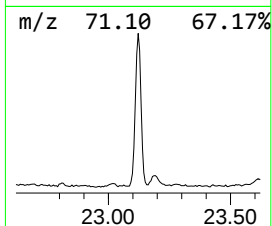
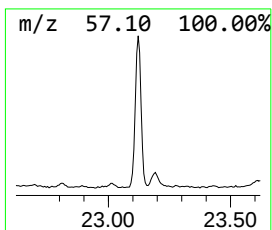
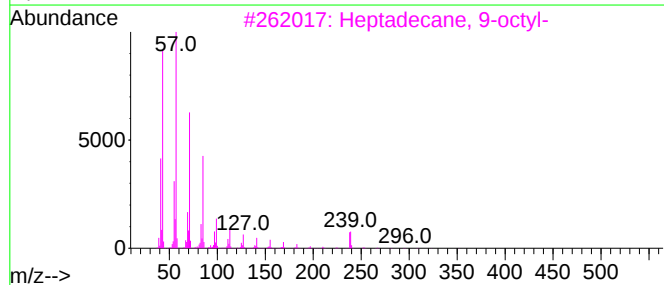
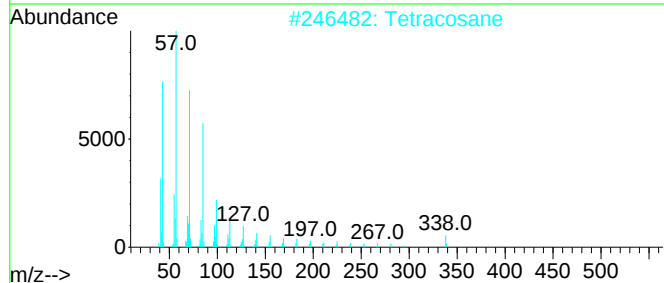
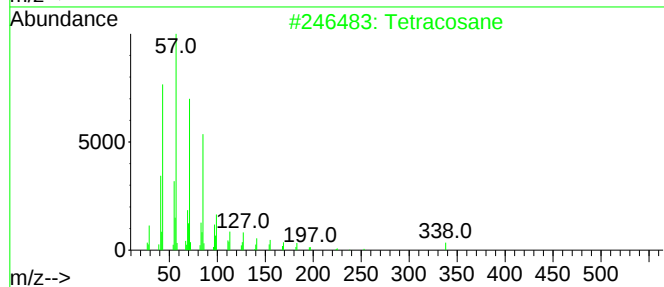
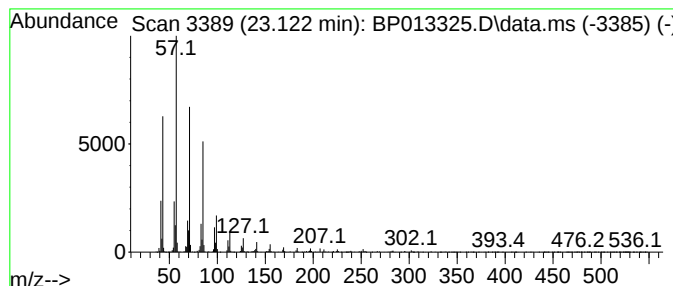
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.122	5.13 ng/ul	1801080	Perylene-d12	23.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	97
2			Tetracosane	338	C24H50	000646-31-1	96
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
4			Heneicosane	296	C21H44	000629-94-7	95
5			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013325.D
Acq On : 25 Dec 2022 05:54
Operator : CG/JU
Sample : N6018-18
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYK3

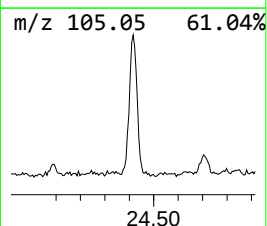
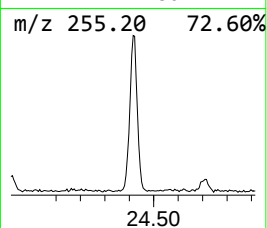
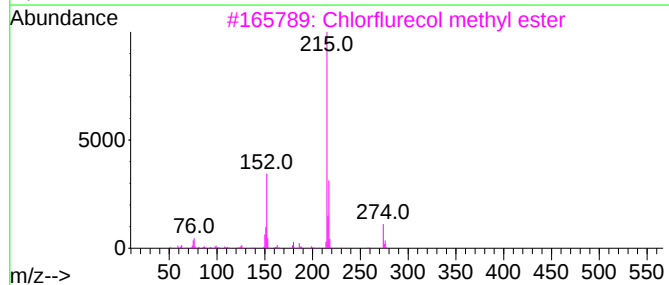
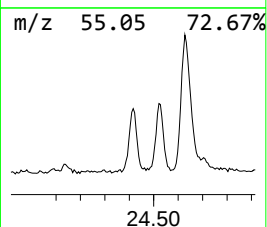
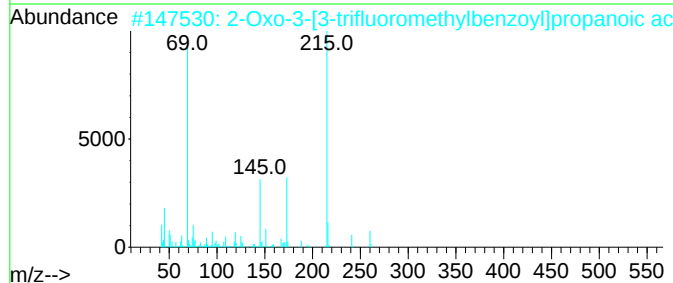
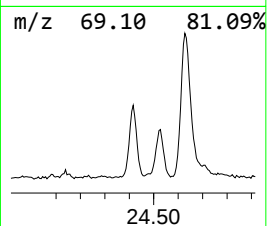
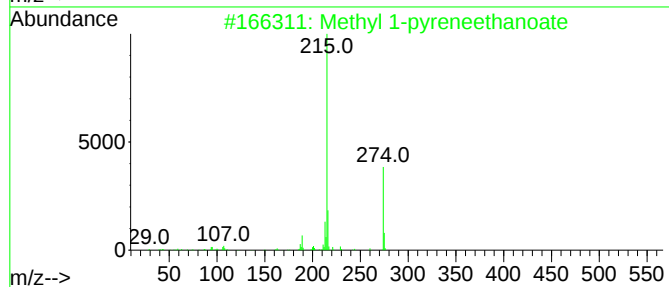
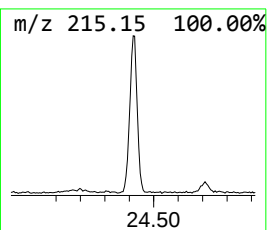
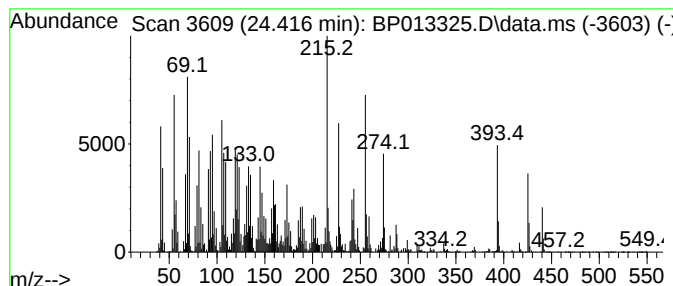
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.416	12.15 ng/ul	4265880	Perylene-d12	23.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	42
2			2-Oxo-3-[3-trifluoromethylbenzoyl]propanoic acid	260	C11H7F3O4	1000211-47-7	38
3			Chlorfluorecol methyl ester	274	C15H11ClO3	002536-31-4	27
4			4-Methoxy-1-naphthyl thiocyanate	215	C12H9NOS	066231-77-4	20
5			2-Methoxy-1-naphthyl thiocyanate	215	C12H9NOS	092696-42-9	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013325.D
Acq On : 25 Dec 2022 05:54
Operator : CG/JU
Sample : N6018-18
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYK3

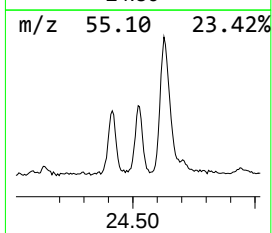
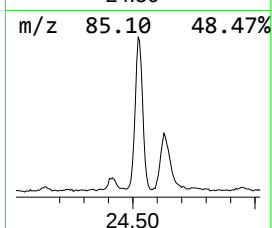
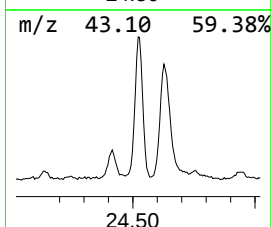
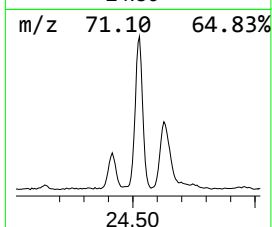
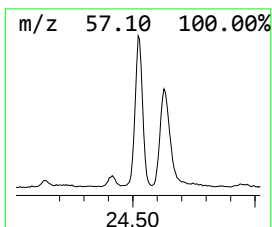
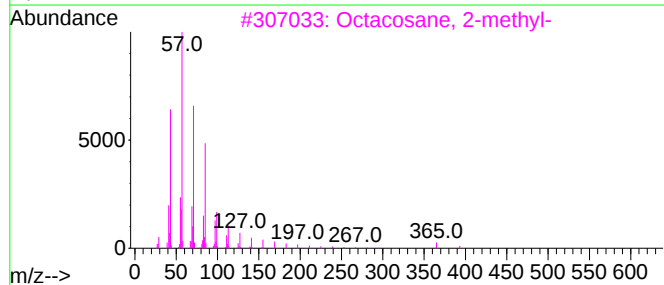
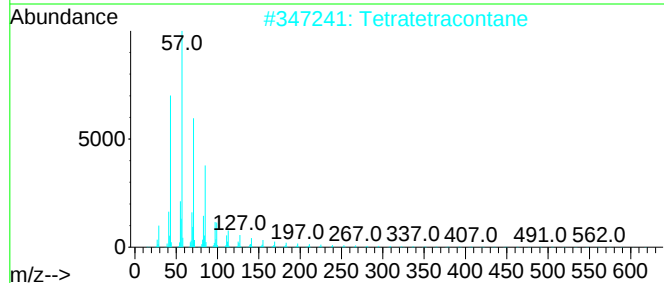
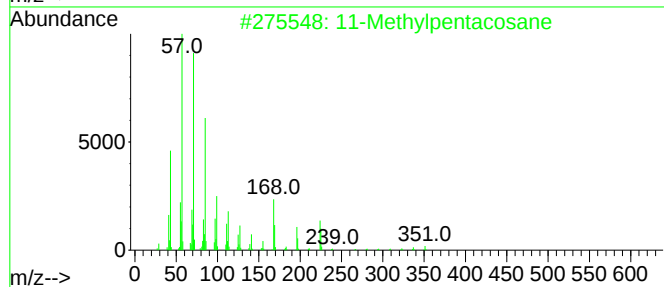
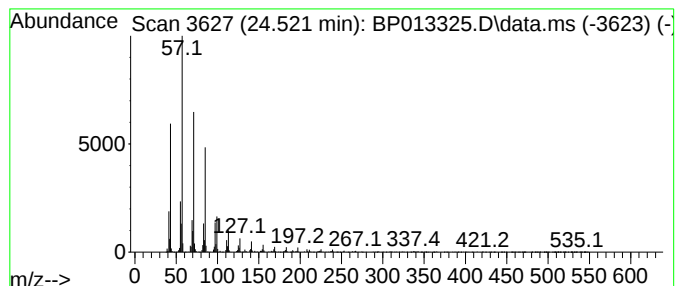
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.522	7.99 ng/ul	2803600	Perylene-d12	23.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11-Methylpentacosane	366	C26H54	015689-71-1	94
2			Tetratetracontane	619	C44H90	007098-22-8	91
3			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91
5			Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013325.D
Acq On : 25 Dec 2022 05:54
Operator : CG/JU
Sample : N6018-18
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYK3

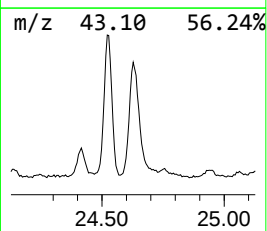
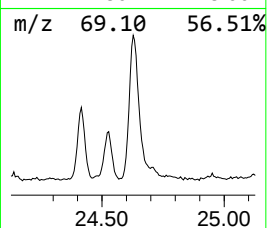
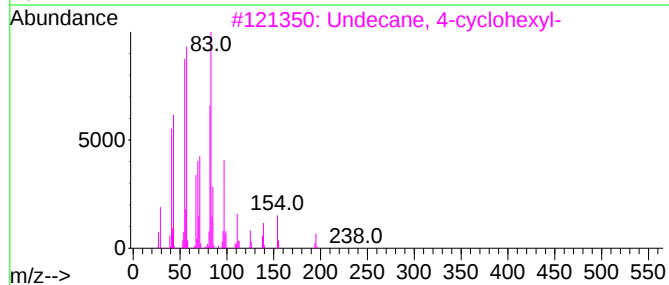
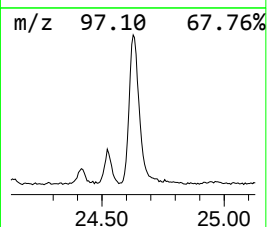
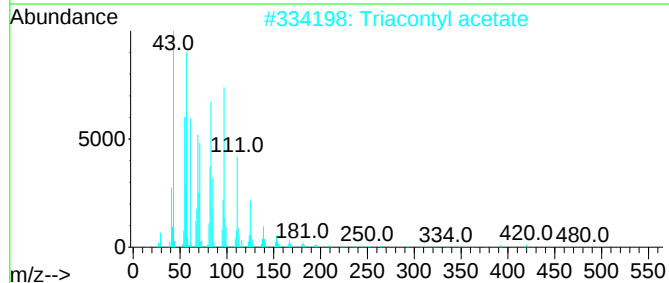
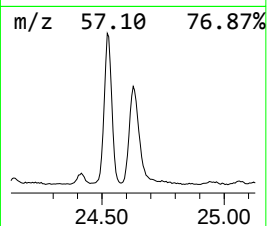
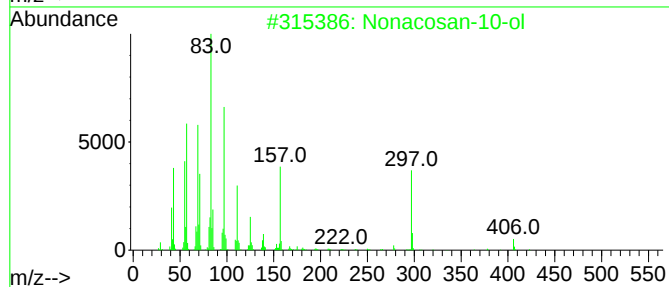
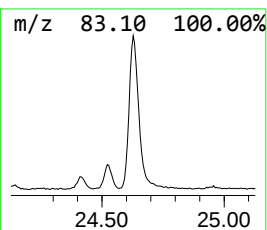
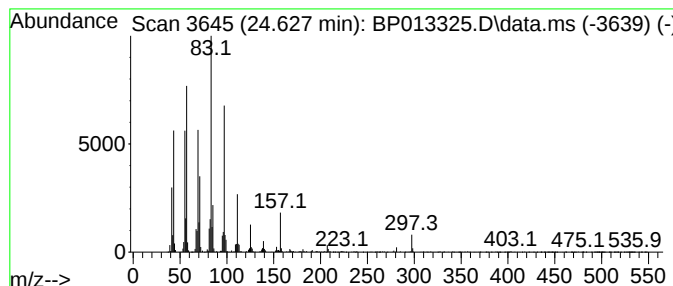
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 Nonacosan-10-ol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.627	12.11 ng/ul	4251110	Perylene-d12	23.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonacosan-10-ol	424	C29H60O	000504-55-2	95
2			Triacontyl acetate	480	C32H64O2	041755-58-2	83
3			Undecane, 4-cyclohexyl-	238	C17H34	013151-79-6	64
4			Tricosyl heptafluorobutyrate	536	C27H47F7O2	1000351-83-4	49
5			Heneicosyl pentafluoropropionate	458	C24H43F5O2	1000351-81-1	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013325.D
Acq On : 25 Dec 2022 05:54
Operator : CG/JU
Sample : N6018-18
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYK3

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
(1R)-2,6,6-Trim...	6.611	32.0	ng/ul	5624370	1	7.940	3519960	20.0
.beta.-Pinene	7.369	7.0	ng/ul	1233200	1	7.940	3519960	20.0
n-Hexadecanoic ...	18.210	5.6	ng/ul	1602520	4	17.340	5693210	20.0
Trichloroacetic...	21.122	11.3	ng/ul	4063710	5	21.433	7173100	20.0
1-Heneicosanol	22.069	2.7	ng/ul	960400	5	21.433	7173100	20.0
(DEL) Alkane: S...	23.122	5.1	ng/ul	1801080	6	23.916	7020690	20.0
unknown-01	24.416	12.2	ng/ul	4265880	6	23.916	7020690	20.0
(DEL) Alkane: S...	24.522	8.0	ng/ul	2803600	6	23.916	7020690	20.0
Nonacosan-10-ol	24.627	12.1	ng/ul	4251110	6	23.916	7020690	20.0