

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
 Data File : BP013322.D
 Acq On : 25 Dec 2022 04:13
 Operator : CG/JU
 Sample : N6018-13
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBYJ3

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: BP013322.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.329	20	24	30	rVB	149671	185592	1.57%	0.136%
2	3.758	94	97	112	rBV	1763145	2546035	21.55%	1.860%
3	4.081	149	152	160	rVB	39660	61033	0.52%	0.045%
4	5.034	308	314	320	rVB2	34360	68471	0.58%	0.050%
5	5.111	325	327	336	rVB10	11055	24225	0.21%	0.018%
6	5.258	348	352	361	rVB	35232	63805	0.54%	0.047%
7	6.452	550	555	561	rVB6	16015	30325	0.26%	0.022%
8	6.528	563	568	575	rBV7	14488	30740	0.26%	0.022%
9	6.611	575	582	589	rVB	1518355	2336009	19.78%	1.706%
10	6.917	629	634	639	rBV4	42802	69247	0.59%	0.051%
11	6.964	639	642	649	rVB7	16633	26687	0.23%	0.019%
12	7.099	656	665	678	rBV	2132159	3432878	29.06%	2.508%
13	7.264	686	693	705	rVB	1789674	2741005	23.20%	2.002%
14	7.369	705	711	718	rBV	187464	295257	2.50%	0.216%
15	7.469	720	728	736	rBV	2266115	3531453	29.90%	2.580%
16	7.940	799	808	815	rBV	1786113	2817195	23.85%	2.058%
17	8.158	839	845	850	rBV	135991	208424	1.76%	0.152%
18	8.211	850	854	859	rVB2	41340	62839	0.53%	0.046%
19	8.434	889	892	897	rVB3	18713	25993	0.22%	0.019%
20	8.652	921	929	940	rBV	2383113	3802251	32.19%	2.777%
21	9.105	999	1006	1013	rBV	2012030	3308705	28.01%	2.417%
22	9.163	1014	1016	1021	rVB3	22062	25729	0.22%	0.019%
23	9.263	1029	1033	1038	rVB4	25806	42455	0.36%	0.031%
24	9.505	1068	1074	1082	rBV3	49153	108874	0.92%	0.080%
25	9.828	1121	1129	1144	rBV	1658981	2783367	23.56%	2.033%
26	10.058	1163	1168	1178	rVB5	53824	115896	0.98%	0.085%
27	10.163	1180	1186	1191	rVB9	13098	24525	0.21%	0.018%
28	10.369	1214	1221	1237	rBV	2416673	4131115	34.97%	3.018%
29	10.752	1277	1286	1292	rBV	2222717	3667036	31.04%	2.679%
30	10.805	1292	1295	1301	rVV	101761	155869	1.32%	0.114%
31	10.893	1301	1310	1330	rVV	1645511	3055164	25.86%	2.232%
32	11.063	1333	1339	1343	rVB9	13433	23321	0.20%	0.017%
33	11.281	1369	1376	1383	rVB9	14652	29741	0.25%	0.022%
34	11.763	1449	1458	1467	rVB9	11998	23272	0.20%	0.017%
35	12.157	1520	1525	1529	rBV2	30964	47878	0.41%	0.035%
36	12.240	1534	1539	1544	rVB5	13706	23402	0.20%	0.017%

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

37	12.334	1550	1555	1560	rBV2	40784	63979	0.54%	0.047%
38	12.410	1564	1568	1578	rVB8	15242	28703	0.24%	0.021%
39	12.769	1626	1629	1637	rVB10	24750	39813	0.34%	0.029%
40	13.110	1681	1687	1690	rVB5	14507	23773	0.20%	0.017%
41	13.340	1718	1726	1730	rBV5	18215	39510	0.33%	0.029%
42	13.393	1731	1735	1741	rVB	57807	81669	0.69%	0.060%
43	13.451	1741	1745	1750	rVB5	37488	57383	0.49%	0.042%
44	13.510	1750	1755	1757	rBV4	15743	23246	0.20%	0.017%
45	13.899	1816	1821	1826	rBV3	157744	234802	1.99%	0.172%
46	13.987	1828	1836	1844	rVV	3448843	4858791	41.13%	3.549%
47	14.051	1844	1847	1852	rVV5	39815	63532	0.54%	0.046%
48	14.104	1852	1856	1860	rVB6	27445	40708	0.34%	0.030%
49	14.275	1879	1885	1903	rVB	4041885	6012716	50.90%	4.392%
50	14.416	1906	1909	1915	rBV6	35316	61384	0.52%	0.045%
51	14.469	1915	1918	1923	rVB2	33597	44102	0.37%	0.032%
52	14.540	1926	1930	1931	rBV4	20263	25152	0.21%	0.018%
53	14.581	1931	1937	1944	rVB	2676350	3833121	32.45%	2.800%
54	14.646	1944	1948	1953	rBV2	92137	138273	1.17%	0.101%
55	14.787	1966	1972	1977	rBV	1230692	1939626	16.42%	1.417%
56	14.822	1977	1978	1988	rVB3	245691	333273	2.82%	0.243%
57	14.940	1993	1998	2002	rVV2	58636	105334	0.89%	0.077%
58	14.987	2002	2006	2012	rVV6	49984	99687	0.84%	0.073%
59	15.040	2012	2015	2021	rVV7	29209	49055	0.42%	0.036%
60	15.116	2024	2028	2032	rVB6	28466	40561	0.34%	0.030%
61	15.310	2057	2061	2066	rVB5	14816	23087	0.20%	0.017%
62	15.369	2066	2071	2075	rBV8	16304	27551	0.23%	0.020%
63	15.493	2084	2092	2099	rBV2	212331	324939	2.75%	0.237%
64	15.575	2099	2106	2121	rBV	4933406	7112020	60.21%	5.195%
65	15.698	2121	2127	2134	rBV	938082	1299178	11.00%	0.949%
66	15.822	2144	2148	2153	rVB4	42188	59919	0.51%	0.044%
67	15.940	2165	2168	2174	rVB	103979	143195	1.21%	0.105%
68	16.022	2178	2182	2184	rBV2	100211	145862	1.23%	0.107%
69	16.151	2198	2204	2210	rBV	239145	336558	2.85%	0.246%
70	16.210	2210	2214	2222	rBV5	88021	123542	1.05%	0.090%
71	16.293	2224	2228	2230	rBV5	15417	24674	0.21%	0.018%
72	16.516	2262	2266	2269	rVB4	29040	38696	0.33%	0.028%
73	16.728	2297	2302	2309	rBV5	88886	157092	1.33%	0.115%
74	16.840	2317	2321	2326	rBV	60177	79480	0.67%	0.058%
75	17.216	2383	2385	2389	rBV5	19195	28133	0.24%	0.021%
76	17.340	2400	2406	2416	rVV	2927288	4397783	37.23%	3.212%

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77	17.440	2416	2423	2431	rVB	4731309	6756722	57.20%	4.936%
78	17.692	2463	2466	2469	rBV3	56547	76742	0.65%	0.056%
79	17.998	2516	2518	2522	rVB4	31273	29650	0.25%	0.022%
80	18.104	2531	2536	2541	rBV8	35567	75477	0.64%	0.055%
81	18.157	2542	2545	2550	rBV	51867	69342	0.59%	0.051%
82	18.216	2550	2555	2566	rBV	797966	1271921	10.77%	0.929%
83	18.339	2573	2576	2582	rVB6	57689	90337	0.76%	0.066%
84	18.545	2607	2611	2620	rVB2	150005	225782	1.91%	0.165%
85	19.104	2703	2706	2711	rBV4	118370	161327	1.37%	0.118%
86	19.328	2738	2744	2745	rBV2	130510	238326	2.02%	0.174%
87	19.445	2760	2764	2774	rBV	237674	359561	3.04%	0.263%
88	19.675	2797	2803	2819	rBV	7165676	9582880	81.13%	7.000%
89	20.616	2959	2963	2967	rBV	537342	687517	5.82%	0.502%
90	21.128	3044	3050	3053	rBV2	2103118	3956026	33.49%	2.890%
91	21.433	3096	3102	3106	rBV	4486932	6882038	58.26%	5.027%
92	22.039	3202	3205	3207	rBV	551642	624070	5.28%	0.456%
93	22.810	3333	3336	3342	rVB	588538	806744	6.83%	0.589%
94	23.127	3385	3390	3395	rBV	1382989	2151461	18.21%	1.572%
95	23.192	3396	3401	3411	rVB	1334031	2768890	23.44%	2.023%
96	23.751	3490	3496	3510	rVB2	5286264	11812411	100.00%	8.629%
97	23.910	3517	3523	3536	rVB2	2830648	6098945	51.63%	4.455%
98	24.139	3558	3562	3570	rVB	1001633	1799839	15.24%	1.315%
99	24.527	3623	3628	3636	rVB	1235805	2568023	21.74%	1.876%
100	24.633	3640	3646	3662	rVB	1907383	5319550	45.03%	3.886%

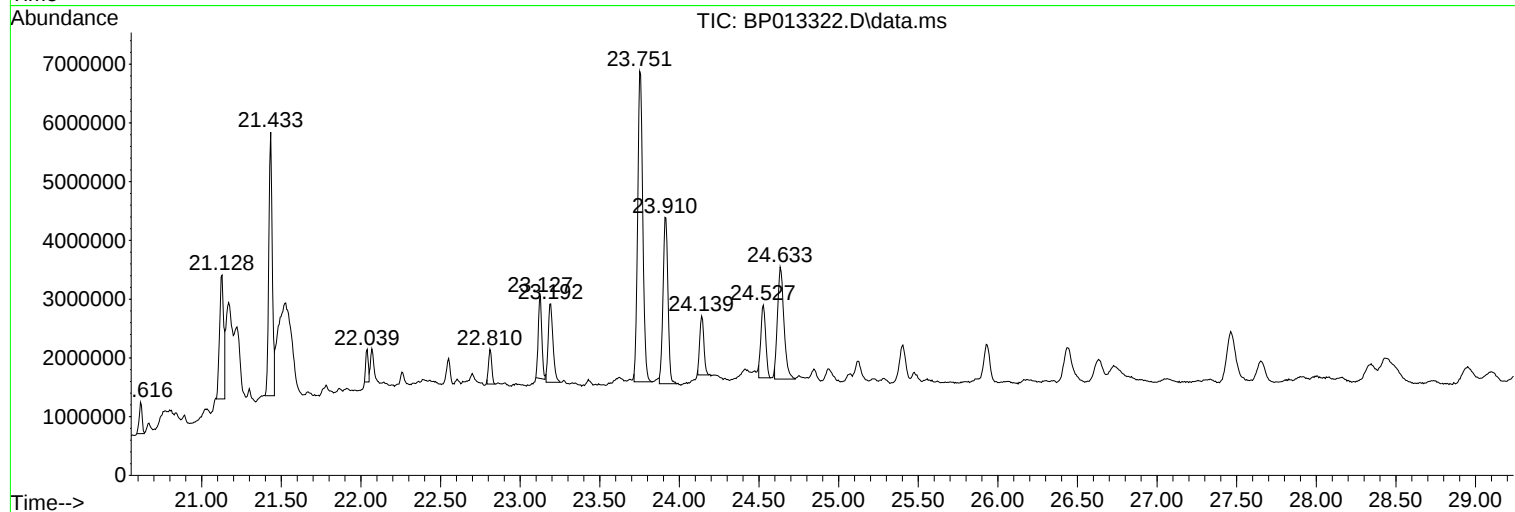
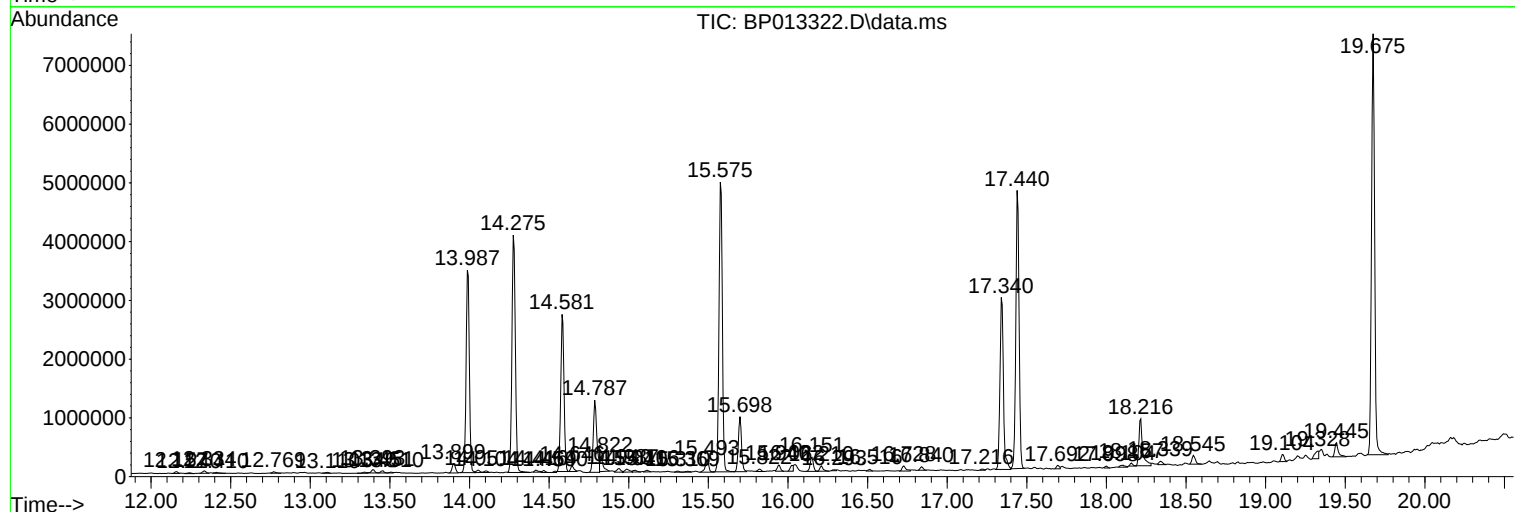
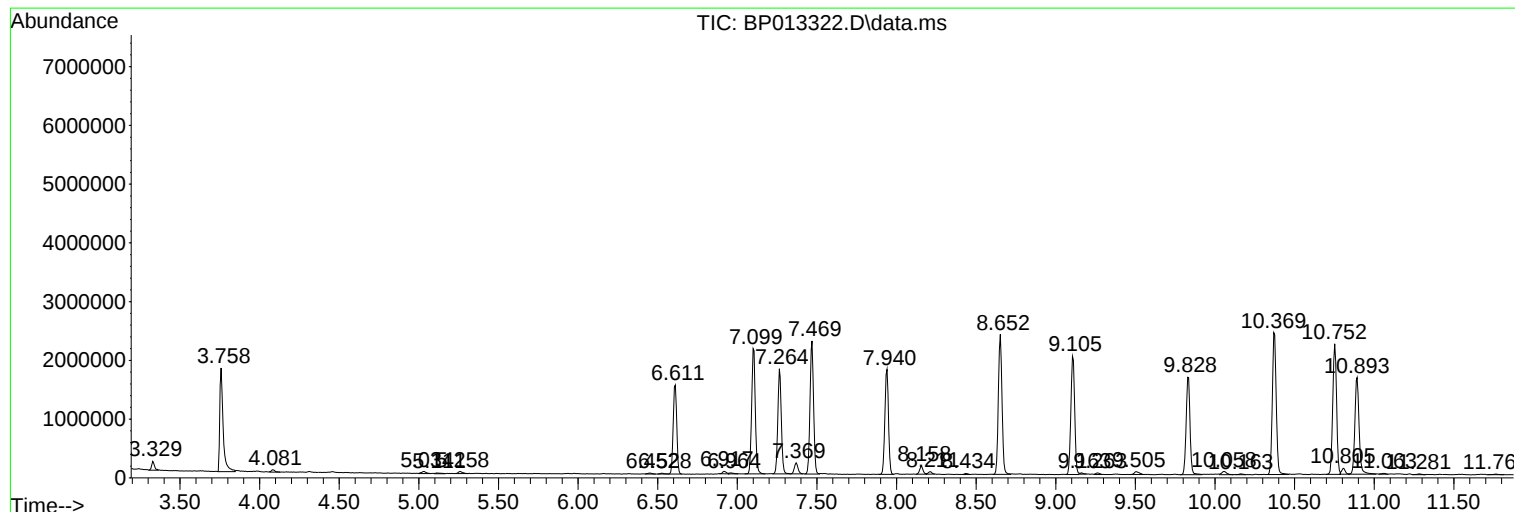
Sum of corrected areas: 136899296

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BNA_P
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DBYJ3

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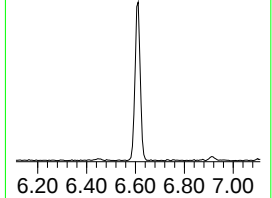
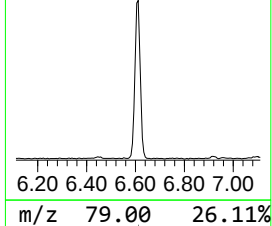
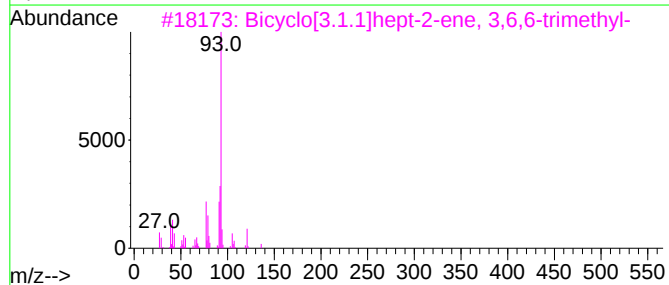
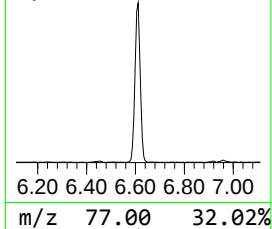
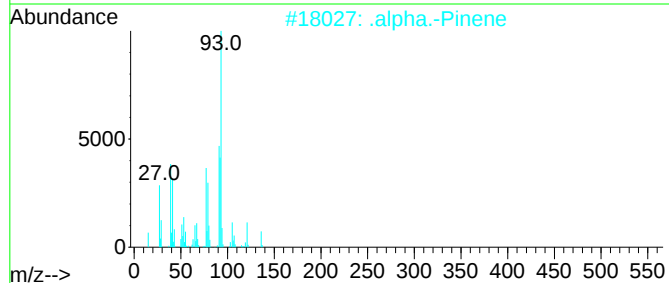
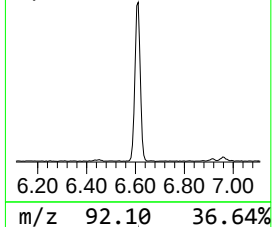
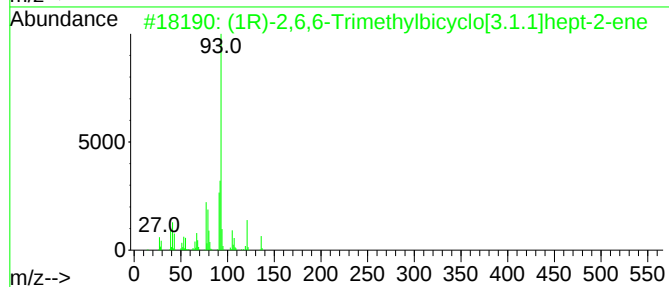
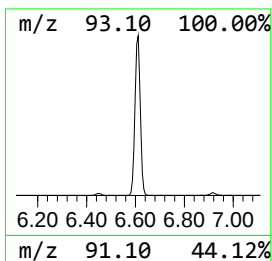
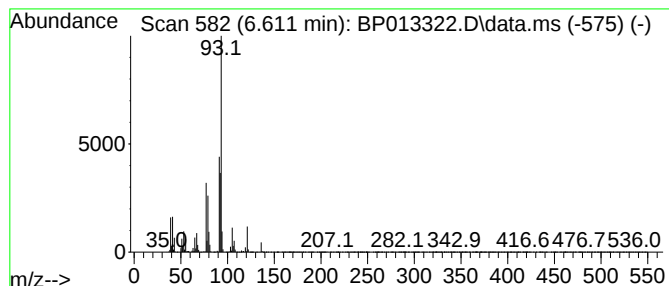
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.611	16.58 ng/ul	2336010	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	95		
3	Bicyclo[3.1.1]hept-2-ene, 3,6,6-trimethyl-	136	C10H16	004889-83-2	94		
4	.alpha.-Pinene	136	C10H16	000080-56-8	94		
5	3-Carene	136	C10H16	013466-78-9	91		



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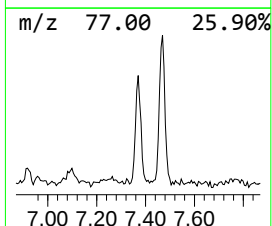
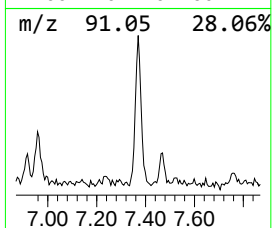
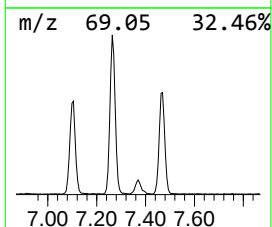
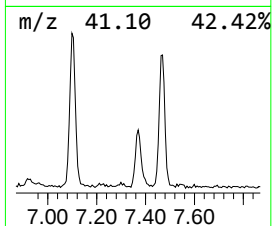
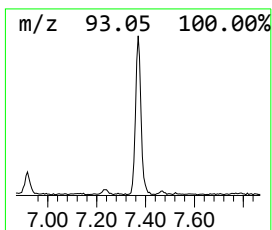
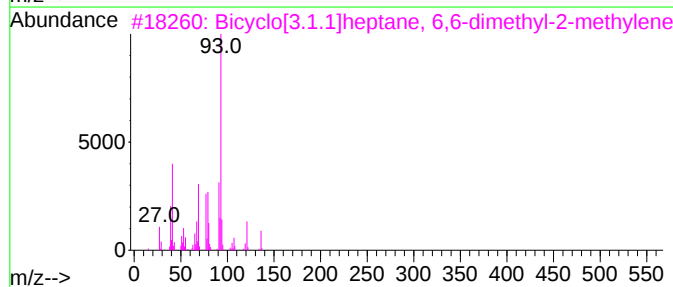
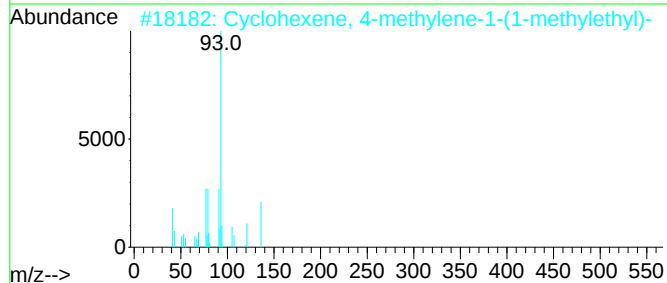
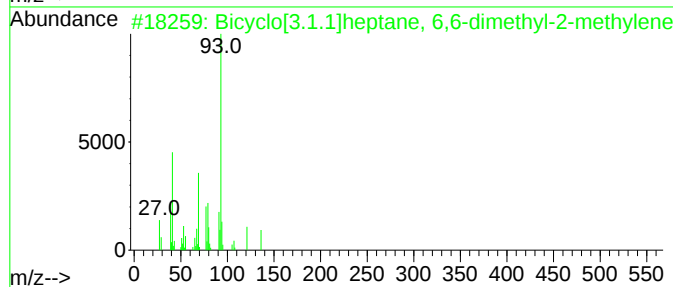
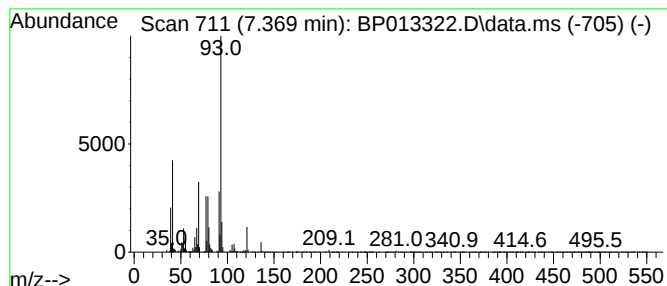
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Bicyclo[3.1.1]heptane, 6,6-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.369	2.10 ng/ul	295257	1,4-Dichlorobenzene-d4	7.940

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



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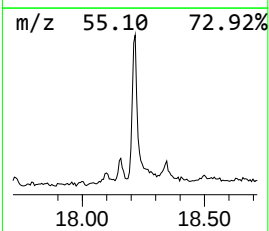
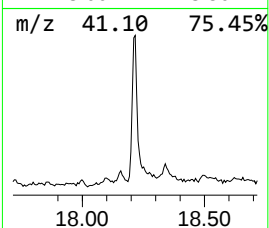
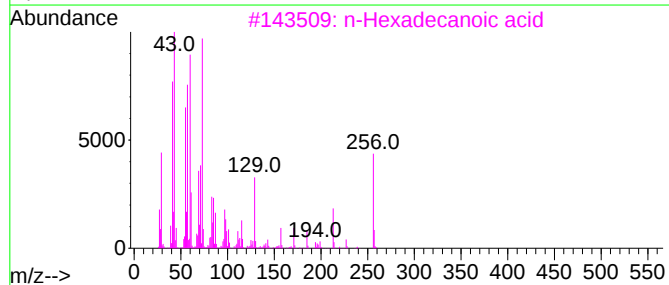
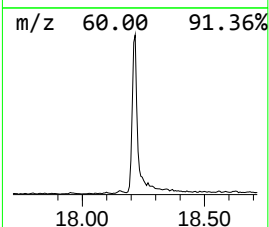
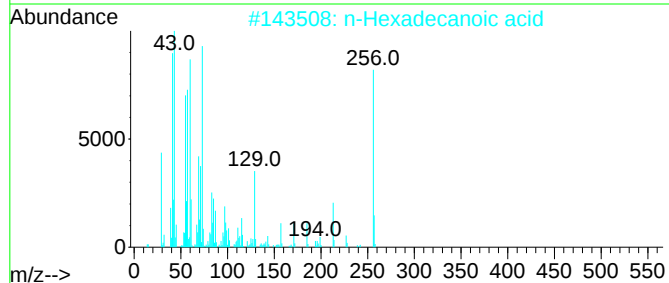
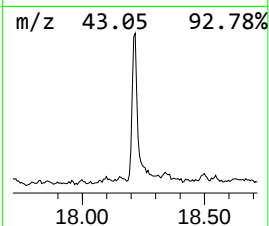
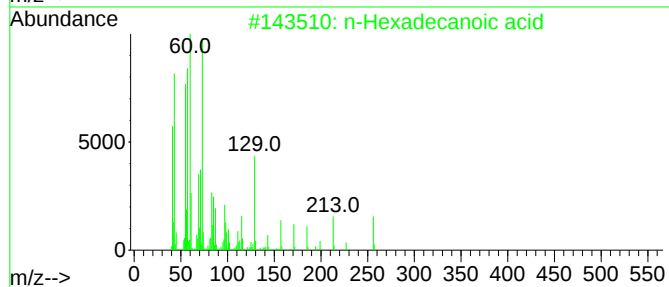
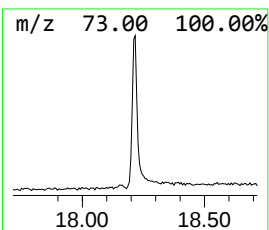
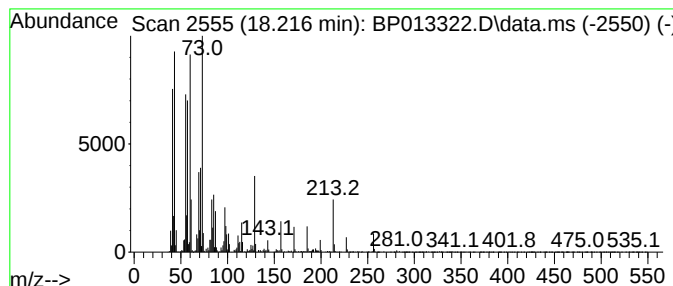
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	5.78 ng/ul	1271920	Phenanthrene-d10	17.340

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
4			Octadecanoic acid	284	C18H36O2	000057-11-4	93
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013322.D
Acq On : 25 Dec 2022 04:13
Operator : CG/JU
Sample : N6018-13
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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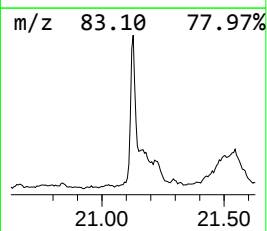
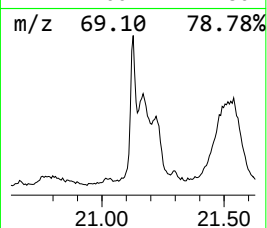
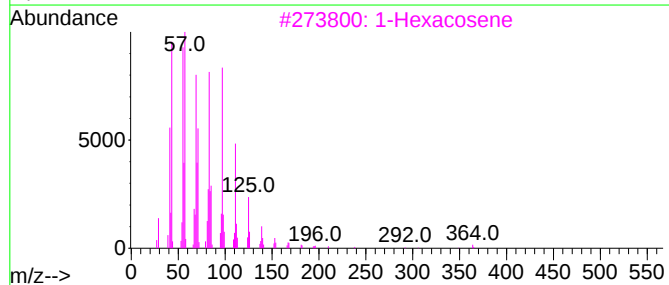
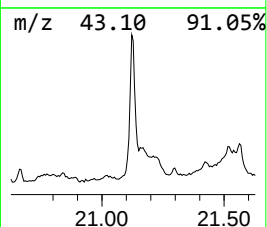
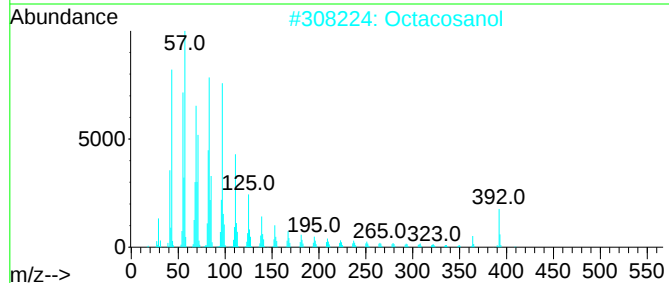
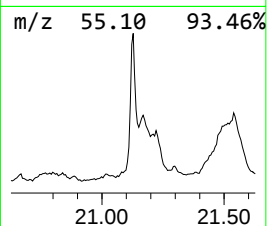
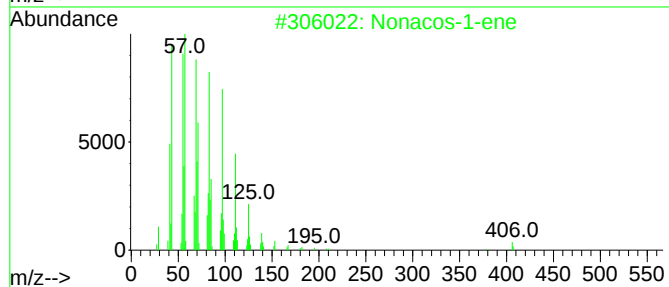
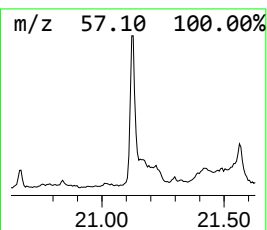
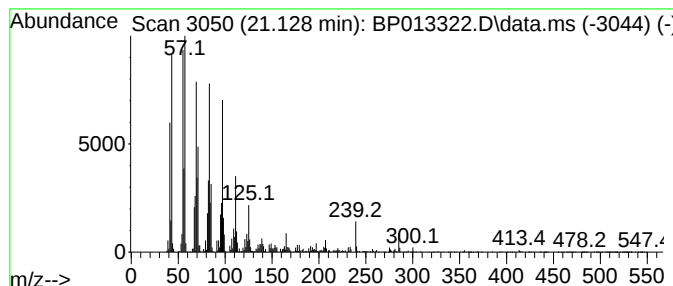
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.128	11.50 ng/ul	3956030	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonacos-1-ene	406	C29H58	018835-35-3	95
2			Octacosanol	410	C28H58O	000557-61-9	94
3			1-Hexacosene	364	C26H52	018835-33-1	94
4			Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	94
5			Pentacos-1-ene	350	C25H50	016980-85-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013322.D
Acq On : 25 Dec 2022 04:13
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Sample : N6018-13
Misc :
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Instrument :
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ClientSampleId :
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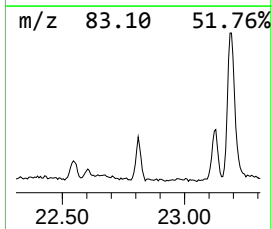
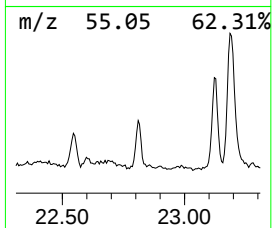
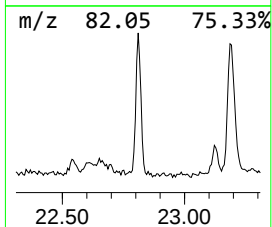
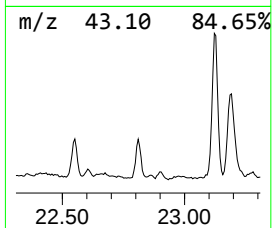
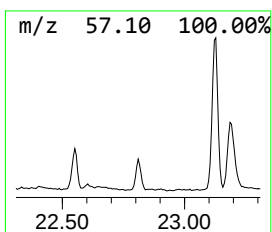
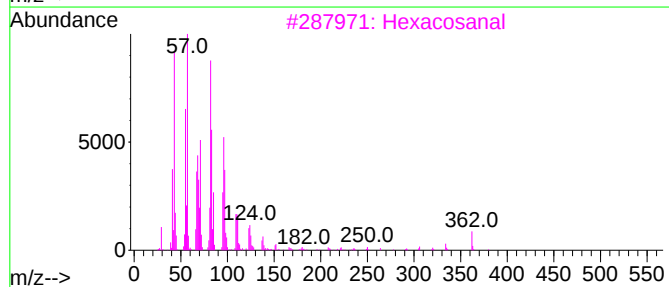
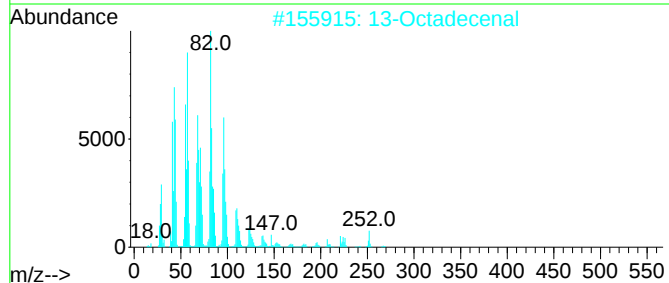
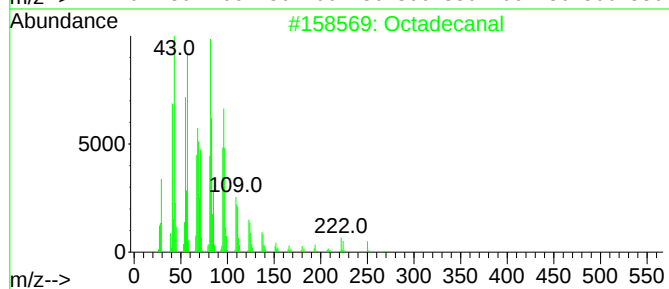
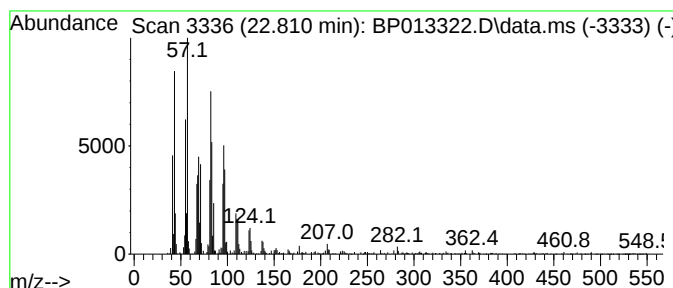
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanal Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.810	2.65 ng/ul	806744	Perylene-d12	23.916

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanal	268	C18H36O	000638-66-4	96
2		13-Octadecenal	266	C18H34O	056554-90-6	94
3		Hexacosanal	380	C26H52O	026627-85-0	94
4		Octadecanal	268	C18H36O	000638-66-4	94
5		Heptadecanal	254	C17H34O	1000376-70-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013322.D
Acq On : 25 Dec 2022 04:13
Operator : CG/JU
Sample : N6018-13
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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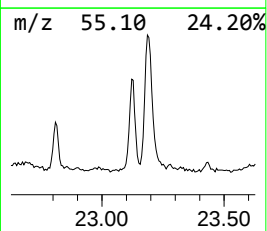
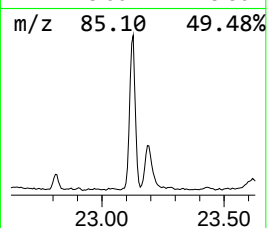
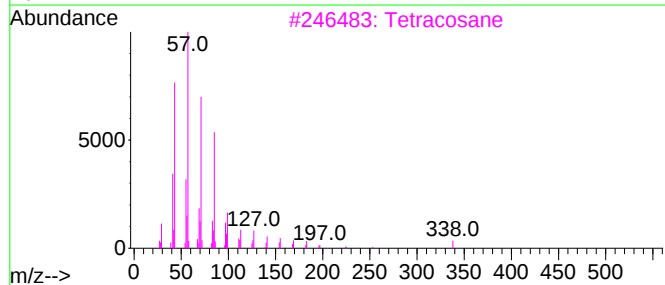
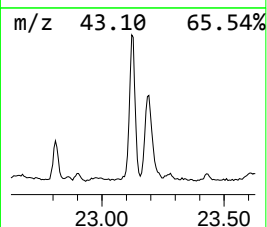
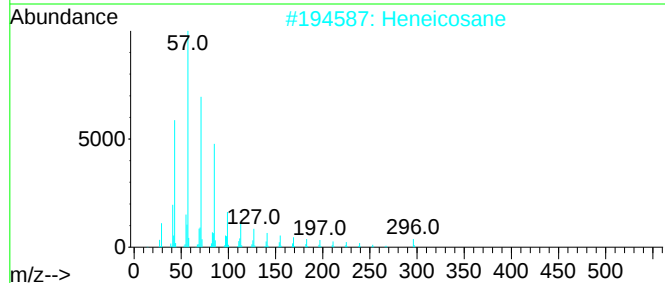
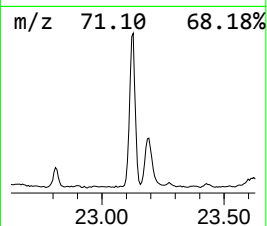
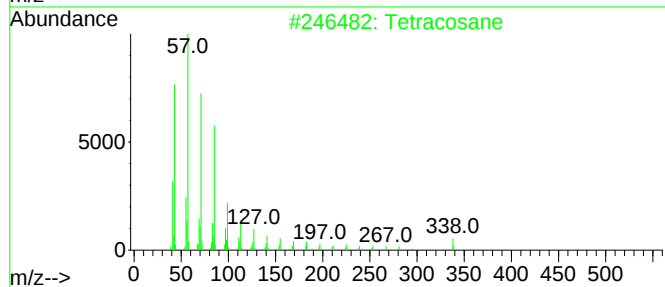
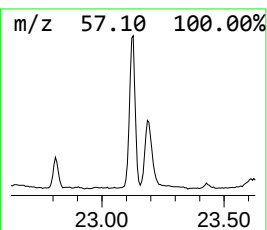
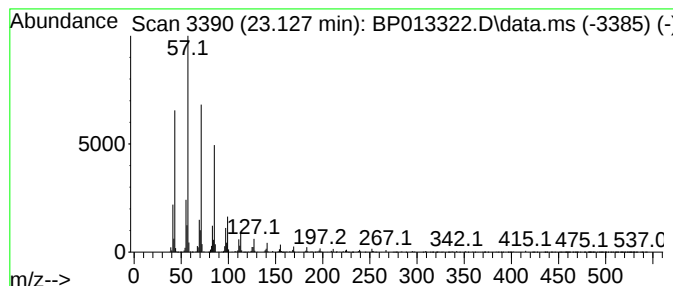
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.128	7.06 ng/ul	2151460	Perylene-d12	23.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	97
2			Heneicosane	296	C21H44	000629-94-7	97
3			Tetracosane	338	C24H50	000646-31-1	96
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
5			Heneicosane, 11-decyl-	437	C31H64	055320-06-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013322.D
Acq On : 25 Dec 2022 04:13
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Instrument :
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ClientSampleId :
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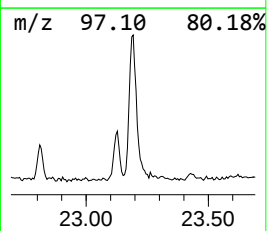
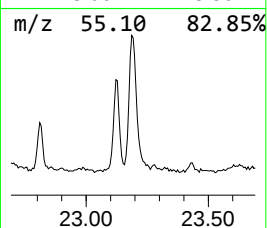
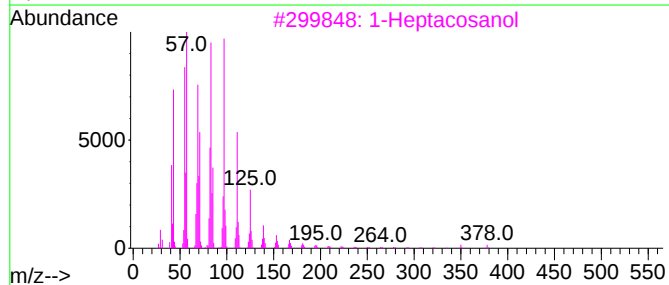
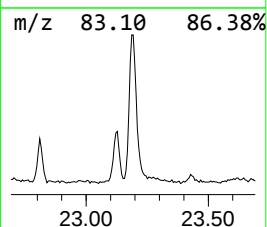
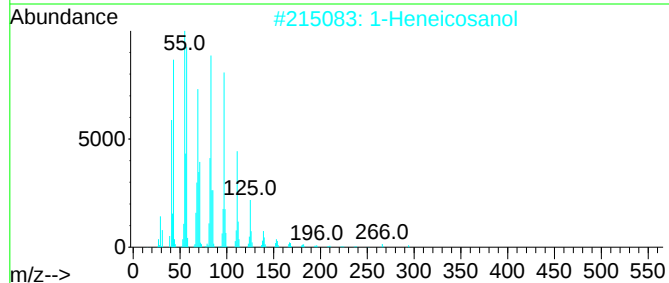
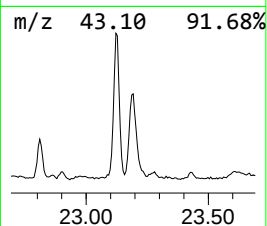
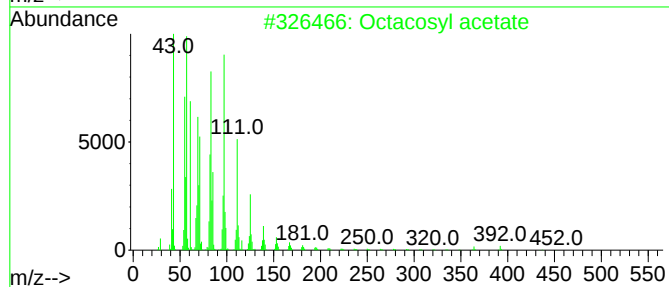
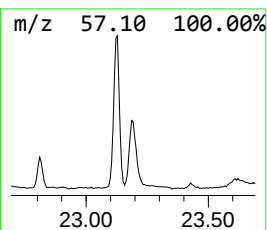
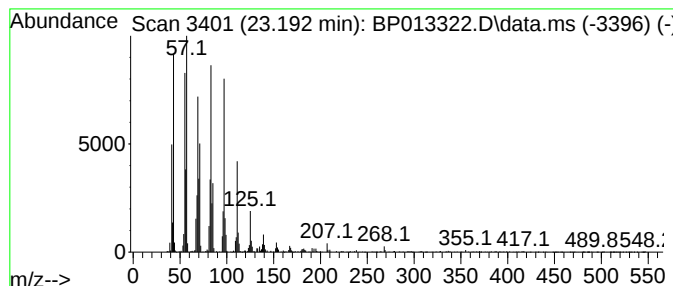
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octacosyl acetate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.192	9.08 ng/ul	2768890	Perylene-d12	23.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosyl acetate	452	C30H60O2	018206-97-8	99
2			1-Heneicosanol	312	C21H44O	015594-90-8	95
3			1-Heptacosanol	396	C27H56O	002004-39-9	94
4			Heptacosyl acetate	438	C29H58O2	1000351-78-2	94
5			Heptacos-1-ene	378	C27H54	015306-27-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
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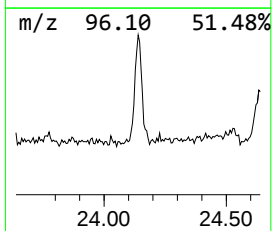
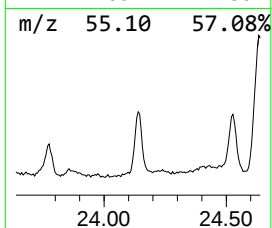
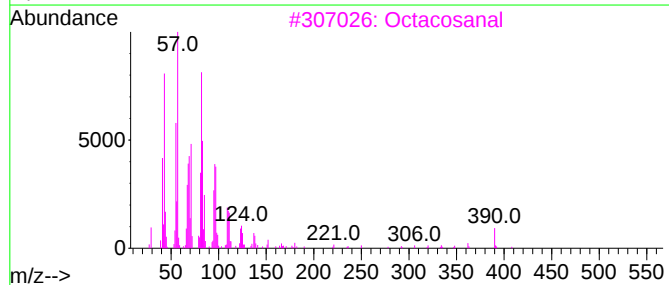
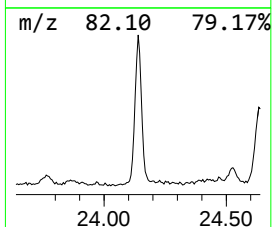
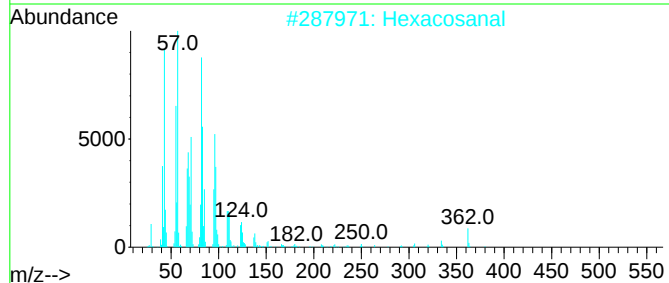
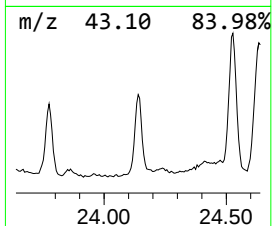
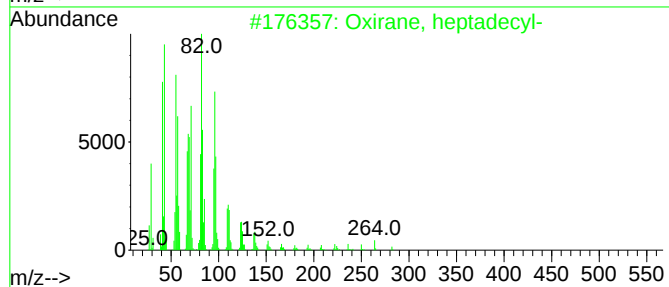
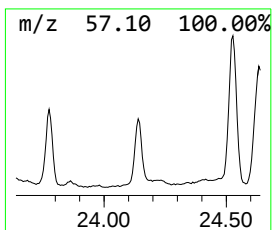
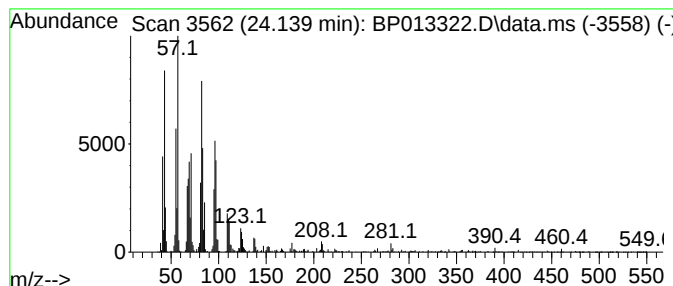
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Oxirane, heptadecyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.139	5.90 ng/ul	1799840	Perylene-d12	23.916	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	95
2	Hexacosanal	380	C26H52O	026627-85-0	94
3	Octacosanal	408	C28H56O	022725-64-0	94
4	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
5	Nonacosanal	422	C29H58O	072934-04-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
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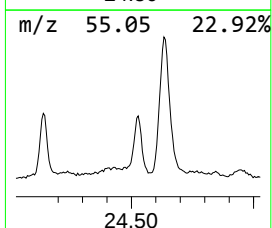
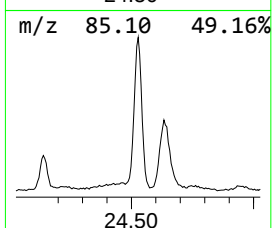
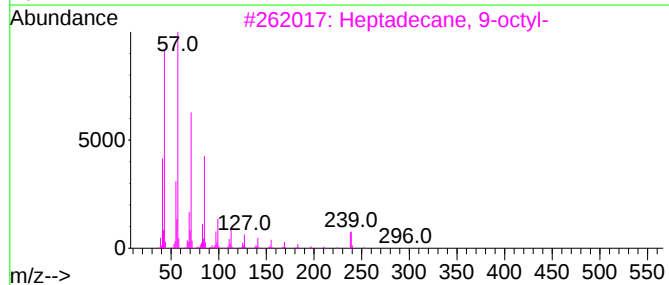
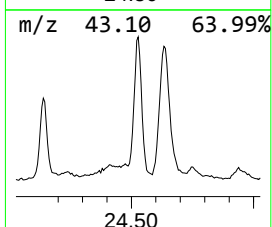
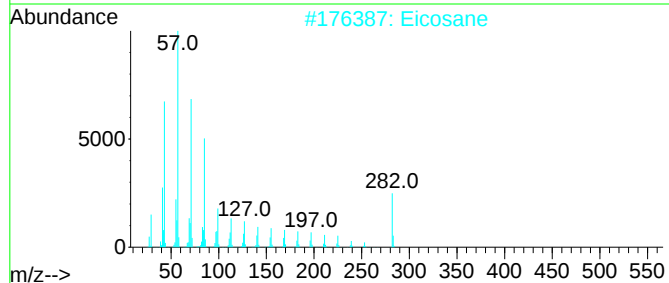
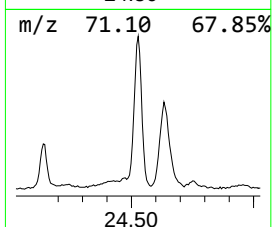
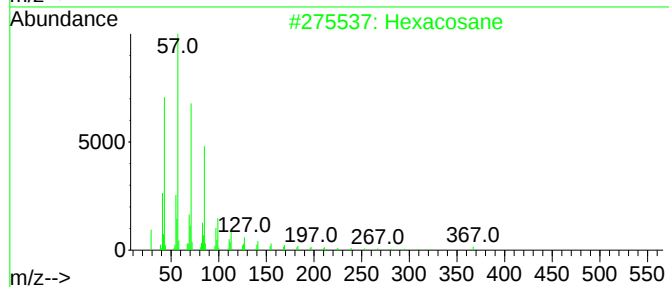
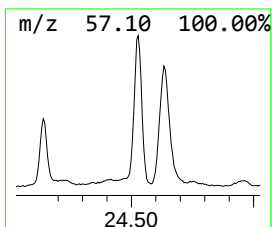
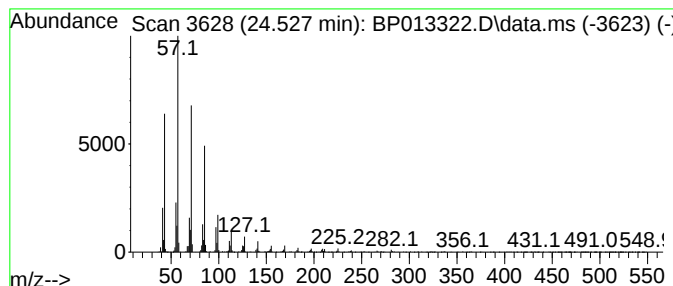
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.527	8.42 ng/ul	2568020	Perylene-d12	23.916

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexacosane	366	C26H54	000630-01-3	94
2		Eicosane	282	C20H42	000112-95-8	93
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
4		Octacosane	394	C28H58	000630-02-4	91
5		Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013322.D
Acq On : 25 Dec 2022 04:13
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Sample : N6018-13
Misc :
ALS Vial : 34 Sample Multiplier: 1

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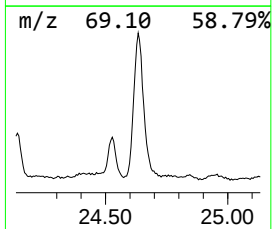
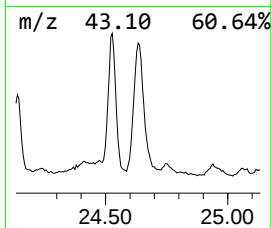
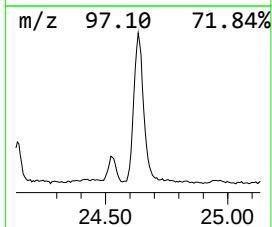
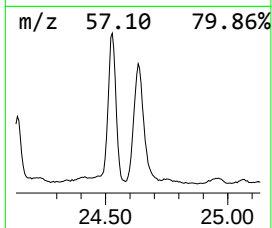
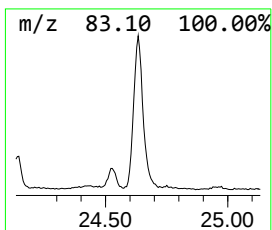
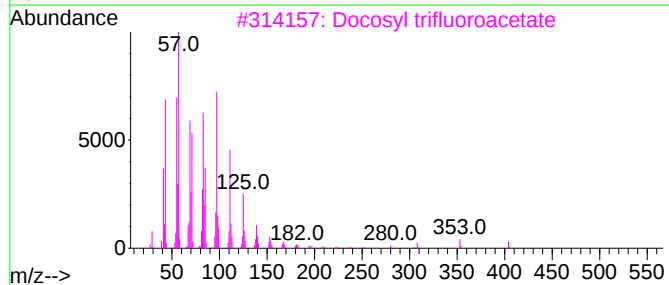
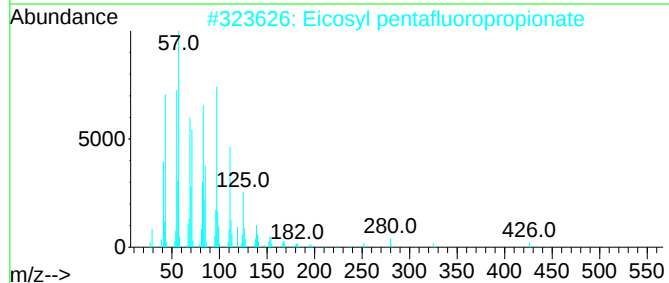
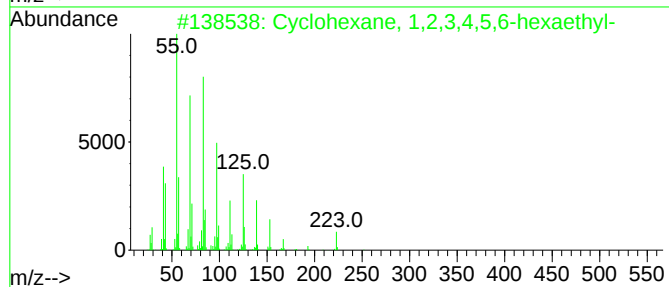
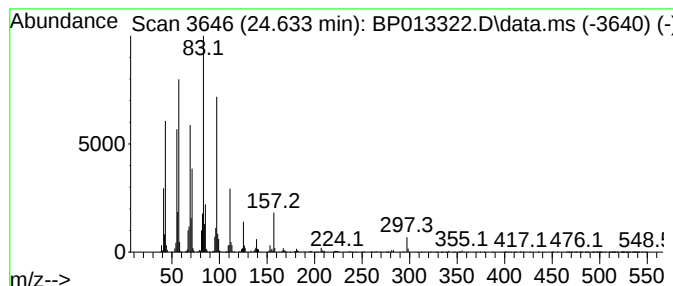
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Cyclic24.633 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.633	17.44 ng/ul	5319550	Perylene-d12	23.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2,3,4,5,6-hexaethyl-	252	C18H36	001795-14-8	64
2			Eicosyl pentafluoropropionate	444	C23H41F5O2	1000351-80-8	58
3			Docosyl trifluoroacetate	422	C24H45F3O2	1000351-75-5	58
4			Pentafluoropropionic acid, octadecyl-	416	C21H37F5O2	959261-25-7	58
5			Tricosyl heptafluorobutyrate	536	C27H47F7O2	1000351-83-4	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
 Data File : BP013322.D
 Acq On : 25 Dec 2022 04:13
 Operator : CG/JU
 Sample : N6018-13
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBYJ3

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	16.6	ng/ul	2336010	1	7.940	2817200	20.0
Bicyclo[3.1.1]h...	7.369	2.1	ng/ul	295257	1	7.940	2817200	20.0
n-Hexadecanoic ...	18.216	5.8	ng/ul	1271920	4	17.340	4397780	20.0
(DEL) Alkane: S...	21.128	11.5	ng/ul	3956030	5	21.433	6882040	20.0
Octadecanal	22.810	2.6	ng/ul	806744	6	23.916	6098950	20.0
(DEL) Alkane: S...	23.128	7.1	ng/ul	2151460	6	23.916	6098950	20.0
Octacosyl acetate	23.192	9.1	ng/ul	2768890	6	23.916	6098950	20.0
Oxirane, heptad...	24.139	5.9	ng/ul	1799840	6	23.916	6098950	20.0
(DEL) Alkane: S...	24.527	8.4	ng/ul	2568020	6	23.916	6098950	20.0
(DEL) Alkane: C...	24.633	17.4	ng/ul	5319550	6	23.916	6098950	20.0