

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013135.D
 Acq On : 19 Dec 2022 20:04
 Operator : CG/JU
 Sample : N6006-13
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXZ7

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: BP013135.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.340	22	26	37	rVB	145467	201895	0.49%	0.099%
2	3.781	98	101	110	rBV	158799	242362	0.59%	0.118%
3	3.875	113	117	122	rVB	67579	89193	0.22%	0.044%
4	3.999	133	138	143	rVB	83301	124911	0.30%	0.061%
5	4.099	151	155	158	rBV	41802	52207	0.13%	0.025%
6	5.040	309	315	326	rBV	770114	1084340	2.64%	0.529%
7	5.505	389	394	401	rVB	102086	151278	0.37%	0.074%
8	6.628	579	585	590	rVB2	51516	83732	0.20%	0.041%
9	7.110	659	667	682	rBV	2326909	3681393	8.95%	1.796%
10	7.281	689	696	710	rVB	1920584	2968447	7.22%	1.448%
11	7.481	723	730	739	rBV	2302327	3630672	8.83%	1.772%
12	7.957	804	811	818	rBV	2114346	3321206	8.08%	1.621%
13	8.663	922	931	942	rBV	2360272	3783567	9.20%	1.846%
14	8.787	946	952	959	rVB	127864	207699	0.51%	0.101%
15	9.122	1001	1009	1023	rBV	2251550	3669413	8.93%	1.790%
16	9.528	1072	1078	1087	rVB	70729	120932	0.29%	0.059%
17	9.846	1125	1132	1143	rBV	1781867	2961488	7.20%	1.445%
18	9.987	1151	1156	1163	rBV2	19027	42843	0.10%	0.021%
19	10.387	1216	1224	1237	rBV	2878921	4780953	11.63%	2.333%
20	10.769	1279	1289	1296	rBV	2956730	4924738	11.98%	2.403%
21	10.904	1305	1312	1330	rVV	1247564	2307105	5.61%	1.126%
22	11.298	1373	1379	1386	rBV6	27788	49677	0.12%	0.024%
23	11.551	1413	1422	1427	rBV6	22226	49033	0.12%	0.024%
24	12.175	1524	1528	1537	rVB	56682	90385	0.22%	0.044%
25	12.293	1537	1548	1554	rBV2	41114	92834	0.23%	0.045%
26	12.351	1554	1558	1566	rVB3	54050	78372	0.19%	0.038%
27	13.222	1698	1706	1712	rBV	102696	163875	0.40%	0.080%
28	13.410	1732	1738	1746	rVB	97852	140511	0.34%	0.069%
29	13.775	1795	1800	1808	rVB4	19410	42901	0.10%	0.021%
30	13.922	1819	1825	1831	rBV8	40104	65590	0.16%	0.032%
31	14.004	1831	1839	1848	rBV	5005355	6918698	16.83%	3.376%
32	14.116	1854	1858	1863	rBV7	29053	45404	0.11%	0.022%
33	14.292	1881	1888	1897	rBV	5843511	8516924	20.72%	4.156%
34	14.481	1916	1920	1929	rVB2	86593	116193	0.28%	0.057%
35	14.598	1929	1940	1947	rBV	4708250	6778880	16.49%	3.308%
36	14.657	1947	1950	1963	rVB9	47682	104927	0.26%	0.051%

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Stop Thrs : 0

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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

37	14.787	1966	1972	1979	rBV	1478910	2312239	5.62%	1.128%
38	14.945	1995	1999	2004	rVV3	38783	54851	0.13%	0.027%
39	15.004	2004	2009	2022	rVV9	51506	164009	0.40%	0.080%
40	15.381	2067	2073	2077	rBV6	19366	41496	0.10%	0.020%
41	15.504	2087	2094	2099	rBV5	72983	110343	0.27%	0.054%
42	15.592	2101	2109	2116	rBV	7491139	10791416	26.25%	5.266%
43	15.645	2116	2118	2123	rVV3	48032	56429	0.14%	0.028%
44	15.704	2123	2128	2142	rVV	2126217	2971811	7.23%	1.450%
45	15.834	2146	2150	2156	rVB3	49711	84908	0.21%	0.041%
46	15.957	2165	2171	2179	rBV	3418528	4457671	10.84%	2.175%
47	16.039	2181	2185	2186	rVV3	35764	49720	0.12%	0.024%
48	16.057	2186	2188	2191	rVV3	48475	75496	0.18%	0.037%
49	16.098	2192	2195	2201	rVB2	74274	115523	0.28%	0.056%
50	16.163	2201	2206	2213	rVB2	85119	124099	0.30%	0.061%
51	16.298	2225	2229	2237	rVB8	31406	75445	0.18%	0.037%
52	16.481	2256	2260	2264	rBV	93610	128266	0.31%	0.063%
53	16.522	2264	2267	2274	rVB	90452	130400	0.32%	0.064%
54	16.616	2277	2283	2287	rBV2	25268	49415	0.12%	0.024%
55	16.739	2300	2304	2307	rBV5	32250	44825	0.11%	0.022%
56	17.057	2355	2358	2362	rVV4	55518	70638	0.17%	0.034%
57	17.233	2384	2388	2395	rBV5	37832	72507	0.18%	0.035%
58	17.351	2402	2408	2413	rBV	5461272	7845930	19.08%	3.828%
59	17.392	2413	2415	2419	rVV	223490	286108	0.70%	0.140%
60	17.451	2419	2425	2438	rVB	7675970	10642782	25.89%	5.193%
61	18.010	2517	2520	2525	rBV5	44307	73625	0.18%	0.036%
62	18.104	2532	2536	2541	rBV4	80002	127983	0.31%	0.062%
63	18.163	2543	2546	2550	rBV2	79219	97116	0.24%	0.047%
64	18.228	2551	2557	2572	rBV	1922048	2911356	7.08%	1.421%
65	18.510	2602	2605	2608	rBV4	36493	47338	0.12%	0.023%
66	18.551	2609	2612	2618	rBV2	73725	113587	0.28%	0.055%
67	18.633	2622	2626	2629	rBV2	184098	231632	0.56%	0.113%
68	19.110	2702	2707	2711	rBV4	136825	232000	0.56%	0.113%
69	19.210	2721	2724	2727	rBV3	49370	69774	0.17%	0.034%
70	19.263	2730	2733	2737	rVB3	130272	166630	0.41%	0.081%
71	19.357	2738	2749	2760	rBV3	825438	2194856	5.34%	1.071%
72	19.457	2761	2766	2780	rVV	794532	1273153	3.10%	0.621%
73	19.680	2799	2804	2816	rVB	8725075	12535976	30.49%	6.117%
74	19.833	2826	2830	2835	rBV	284683	333681	0.81%	0.163%
75	20.180	2884	2889	2894	rBV4	598147	1337300	3.25%	0.653%
76	20.674	2971	2973	2979	rVB2	156744	171348	0.42%	0.084%

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 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
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77	21.133	3046	3051	3062	rBV	1715486	2671441	6.50%	1.304%
78	21.439	3097	3103	3108	rBV	4777983	6575447	15.99%	3.208%
79	22.051	3204	3207	3208	rBV2	224687	267383	0.65%	0.130%
80	22.074	3208	3211	3221	rVB3	547566	945660	2.30%	0.461%
81	23.139	3387	3392	3397	rBV2	1083209	1886265	4.59%	0.920%
82	23.757	3491	3497	3511	rVB2	4257673	9233743	22.46%	4.506%
83	23.921	3518	3525	3545	rVB2	2935676	6297514	15.32%	3.073%
84	24.545	3626	3631	3639	rVB	904610	1853730	4.51%	0.905%
85	26.462	3950	3957	3977	rVB2	1143549	3882606	9.44%	1.894%
86	27.474	4122	4129	4150	rVB3	1505986	5855798	14.24%	2.857%
87	28.527	4294	4308	4328	rVB2	10289055	41113672	100.00%	20.061%

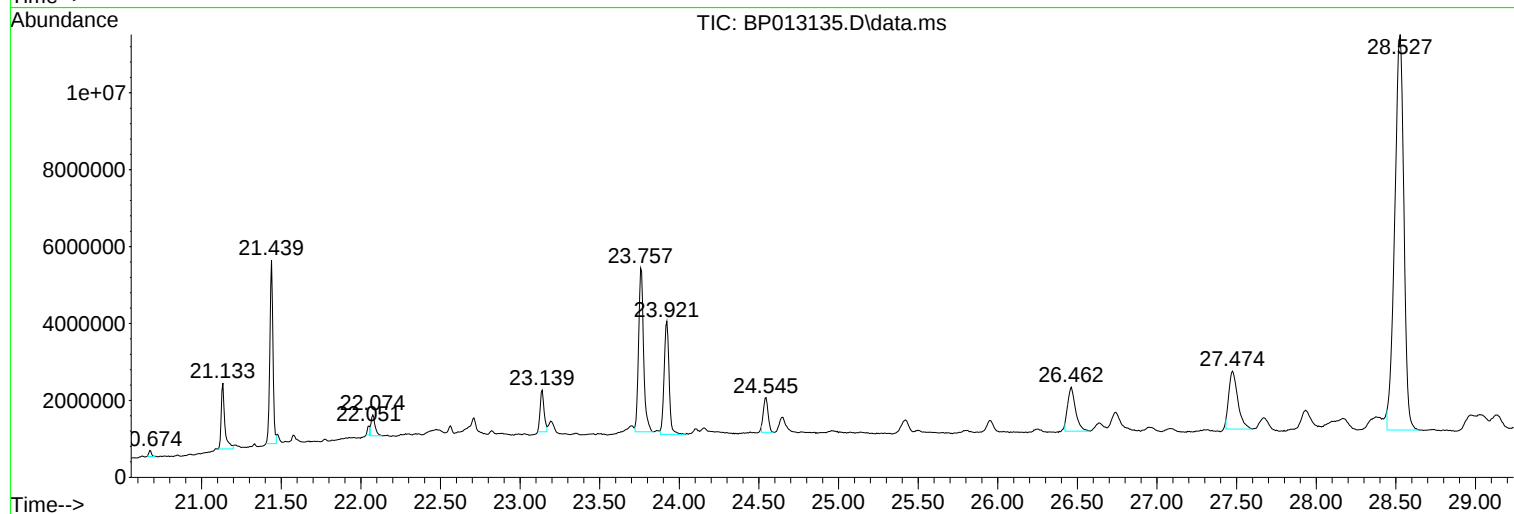
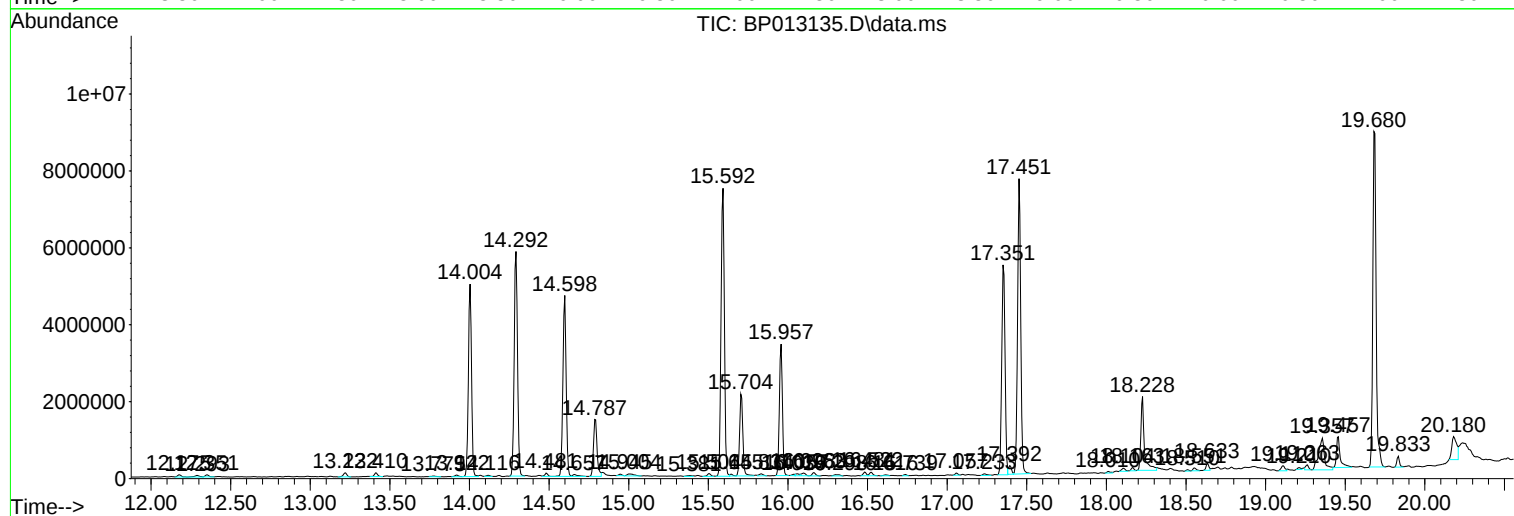
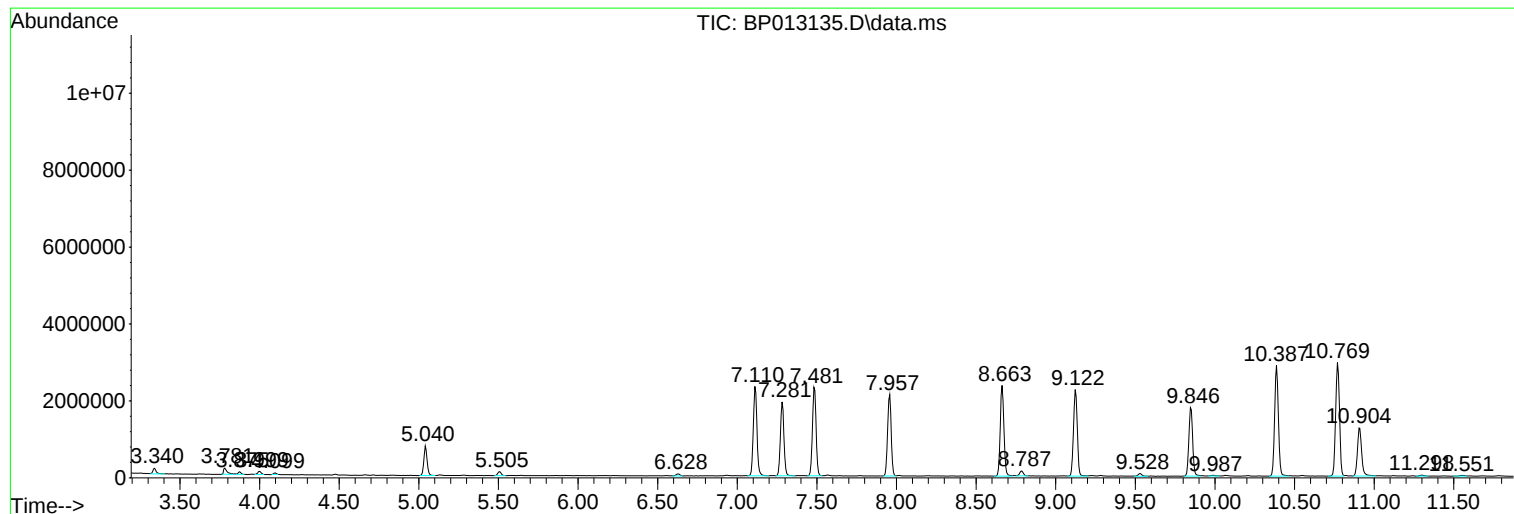
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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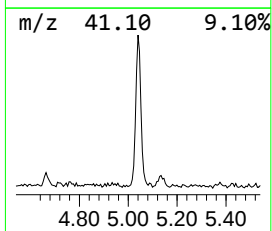
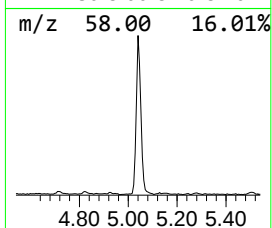
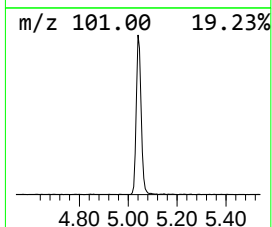
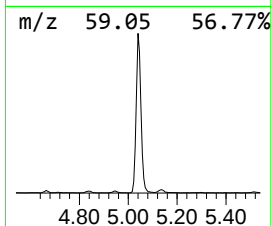
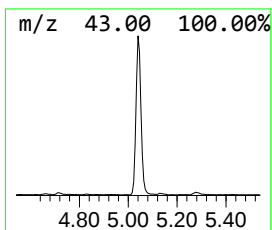
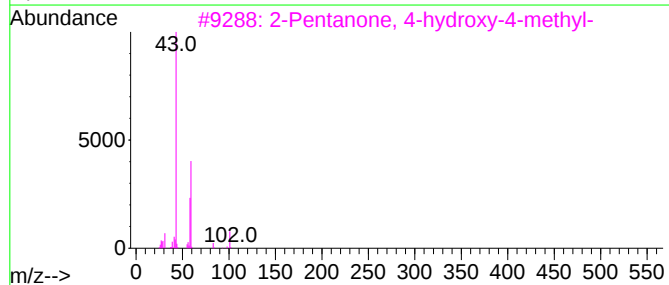
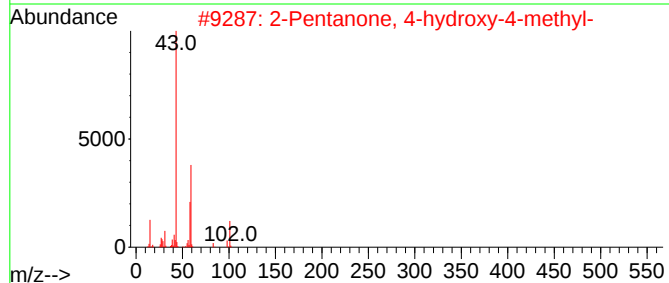
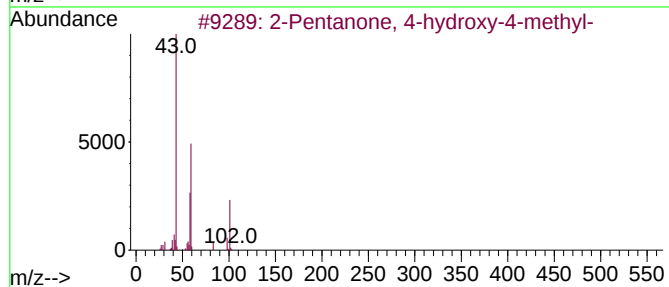
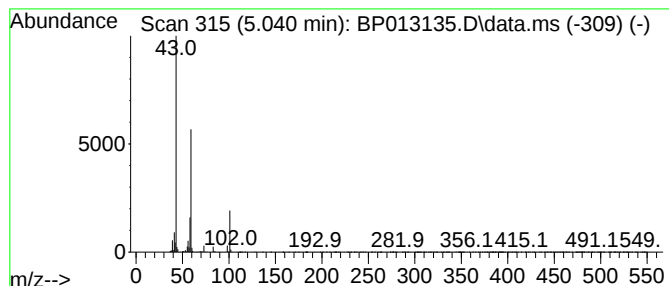
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.040	6.53 ng/ul	1084340	1,4-Dichlorobenzene-d4	7.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		
5	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	25		



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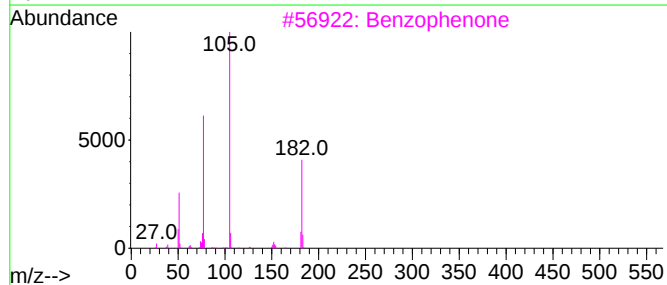
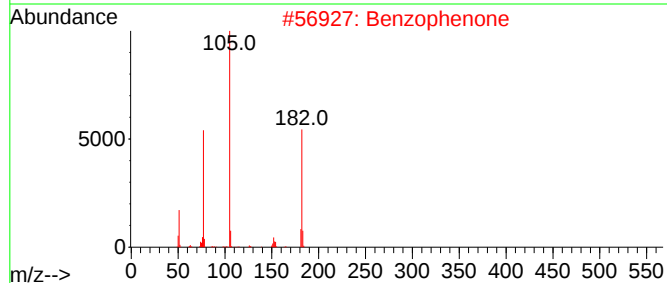
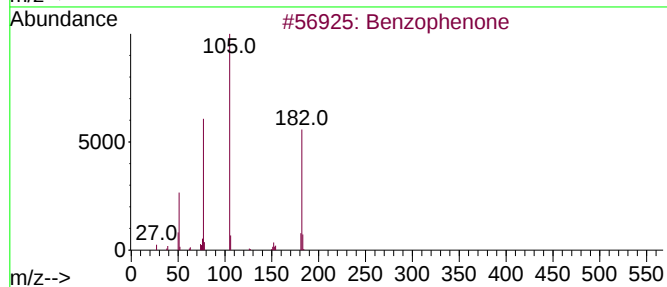
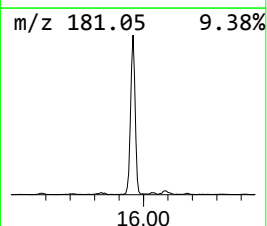
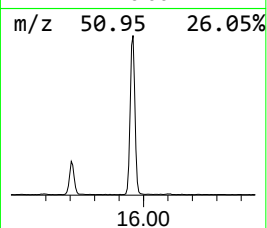
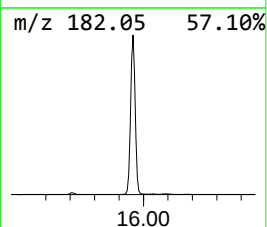
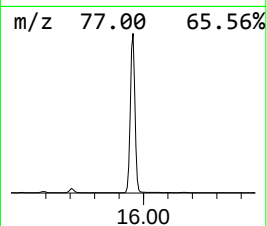
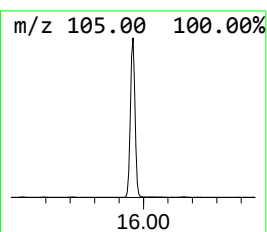
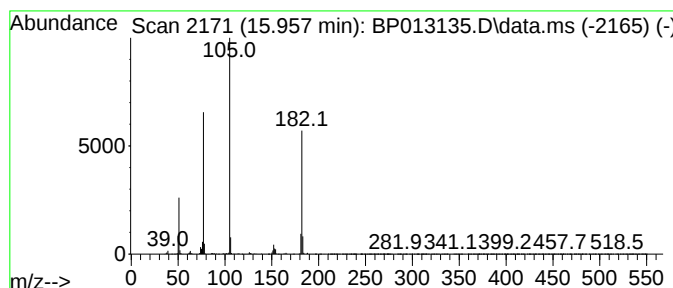
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.957	13.15 ng/ul	4457670	Acenaphthene-d10	14.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	90



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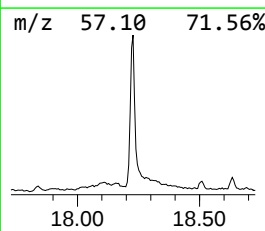
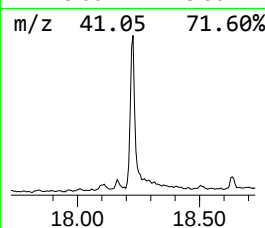
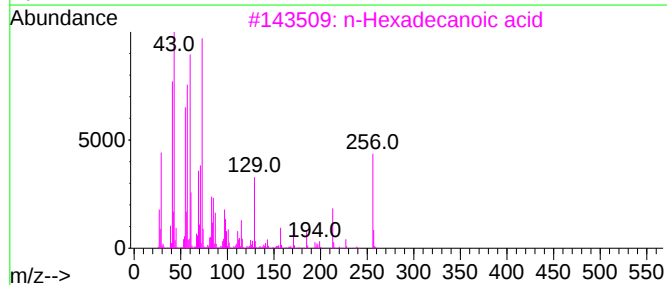
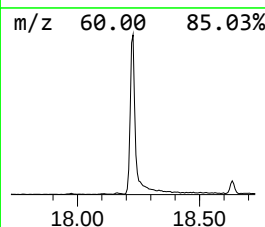
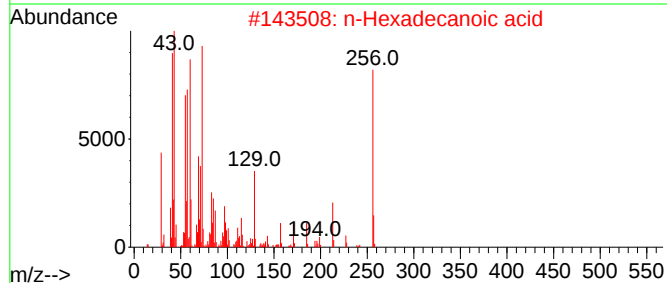
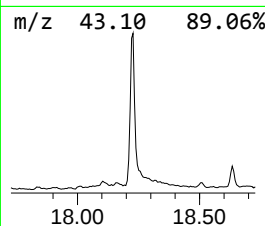
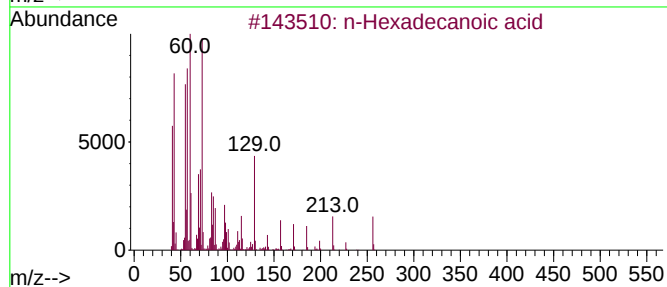
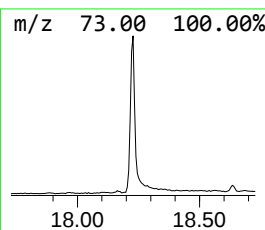
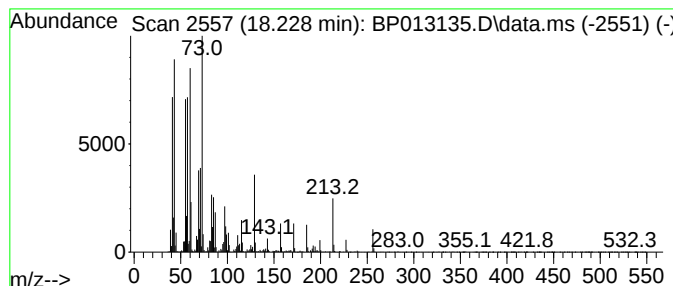
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.227	7.42 ng/ul	2911360	Phenanthrene-d10	17.351

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			Tridecanoic acid	214	C13H26O2	000638-53-9	95
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95



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ClientSampleId :
DBXZ7

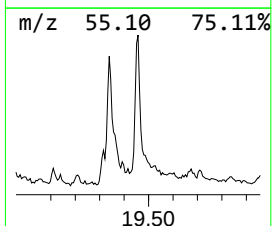
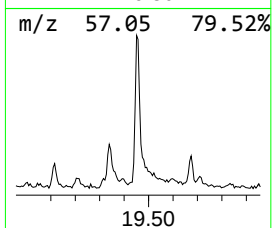
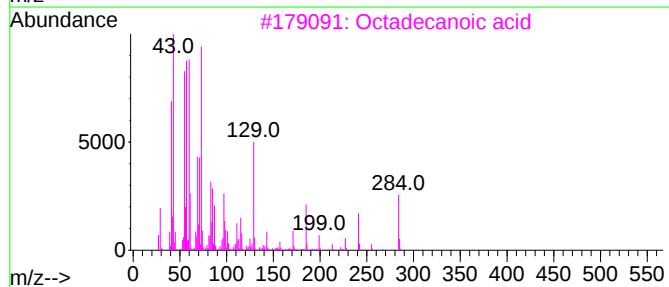
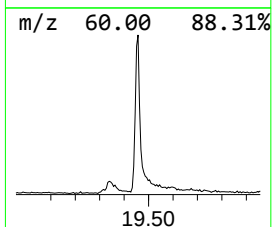
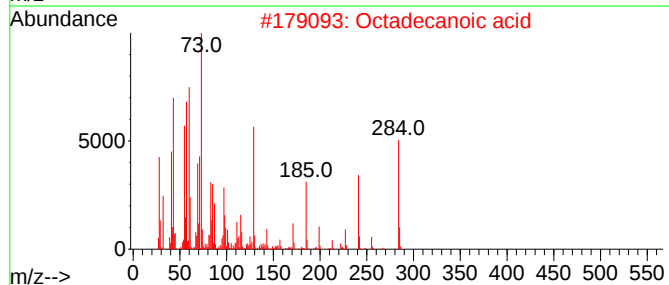
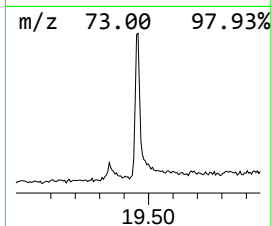
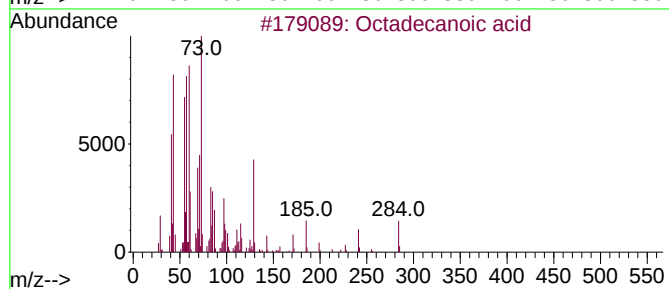
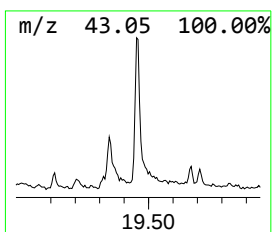
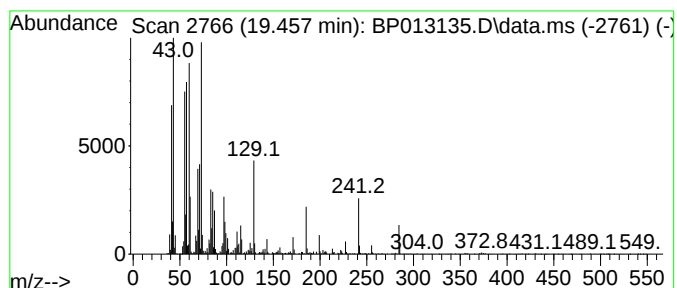
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.457	3.87 ng/ul	1273150	Chrysene-d12	21.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	98
5			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013135.D
Acq On : 19 Dec 2022 20:04
Operator : CG/JU
Sample : N6006-13
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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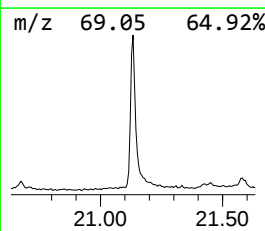
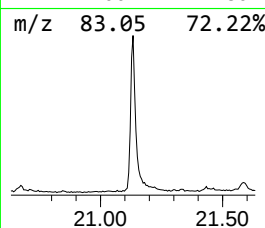
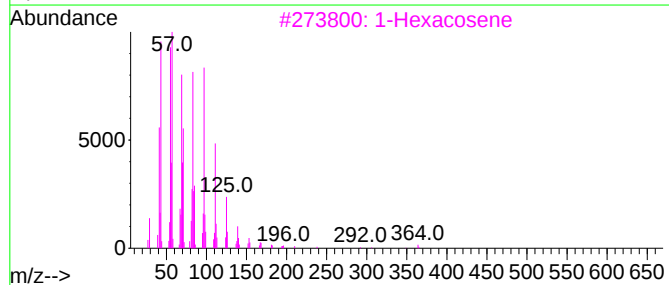
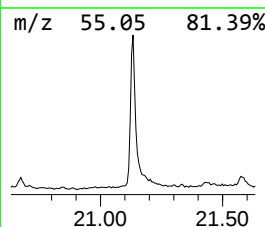
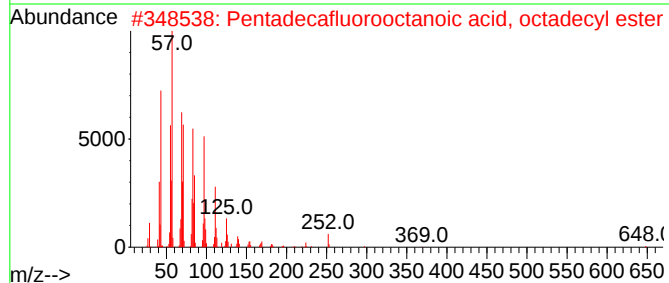
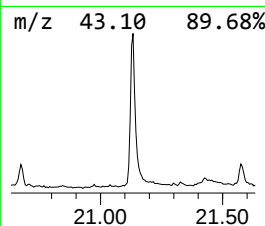
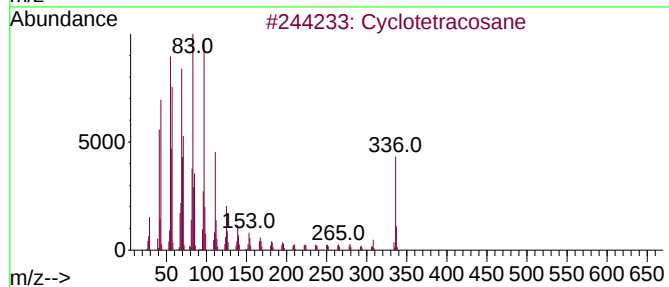
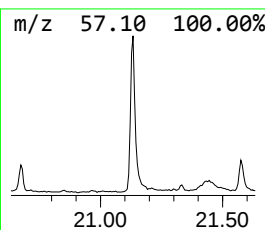
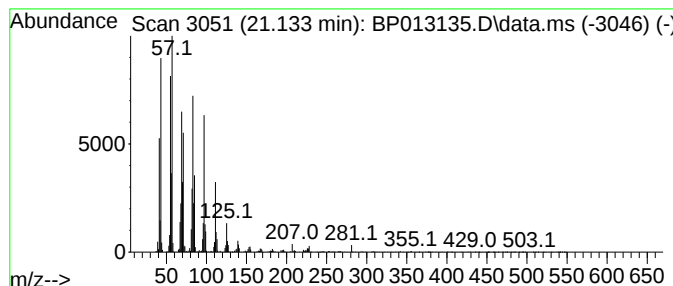
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic21.133 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.133	8.13 ng/ul	2671440	Chrysene-d12	21.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	97
2			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
3			1-Hexacosene	364	C26H52	018835-33-1	94
4			Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	94
5			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013135.D
Acq On : 19 Dec 2022 20:04
Operator : CG/JU
Sample : N6006-13
Misc :
ALS Vial : 14 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

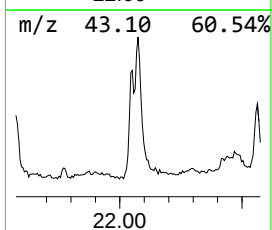
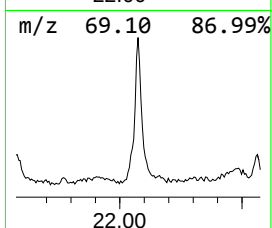
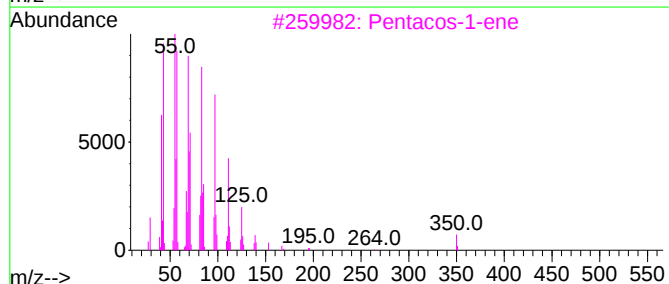
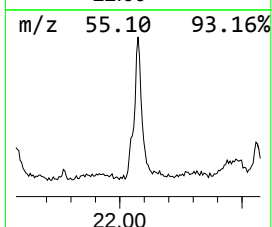
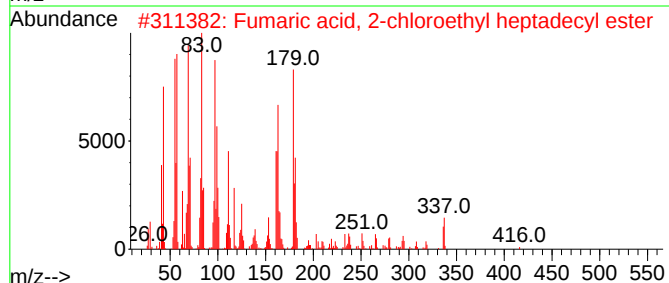
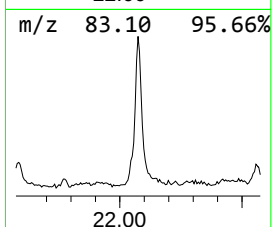
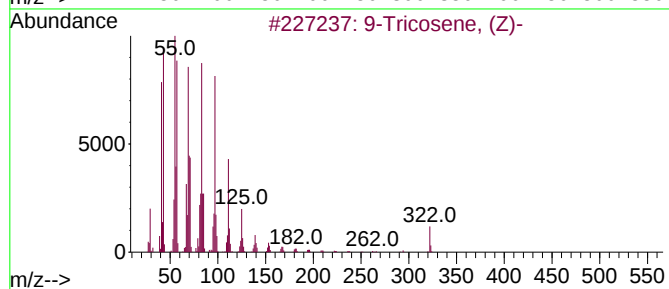
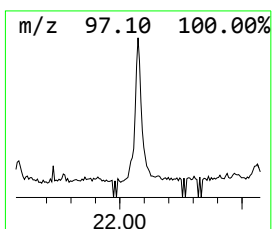
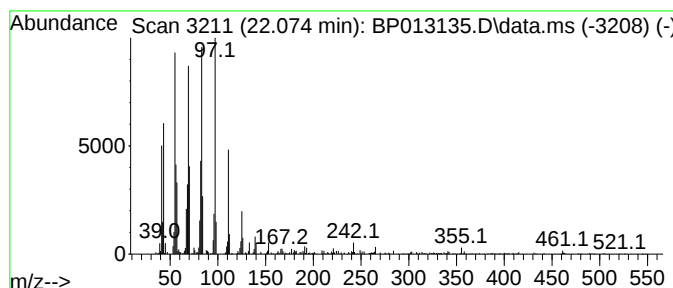
TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.074	2.88 ng/ul	945660	Chrysene-d12	21.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Tricosene, (Z)-			322	C23H46	027519-02-4	91
2	Fumaric acid, 2-chloroethyl hept...			416	C23H41ClO4	1000330-62-3	86
3	Pentacos-1-ene			350	C25H50	016980-85-1	83
4	1-Heneicosanol			312	C21H44O	015594-90-8	76
5	1-Hexadecanol			242	C16H34O	036653-82-4	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013135.D
Acq On : 19 Dec 2022 20:04
Operator : CG/JU
Sample : N6006-13
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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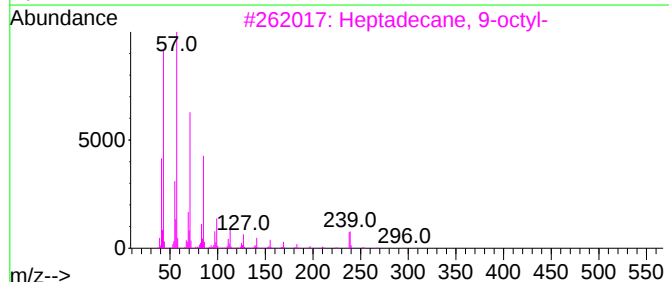
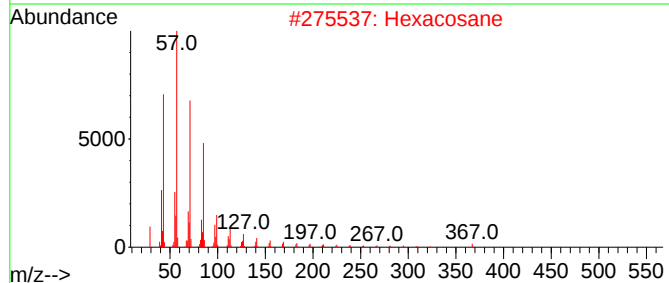
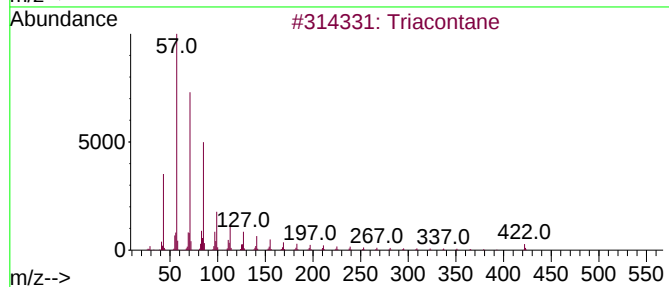
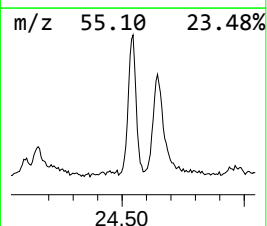
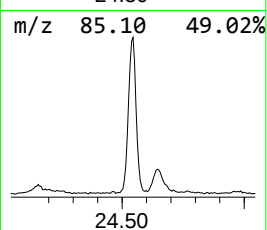
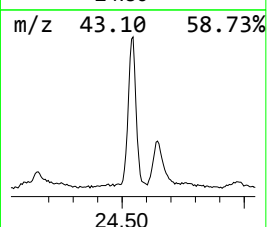
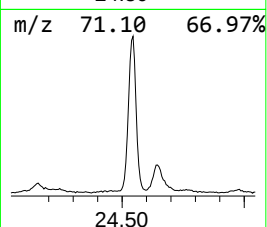
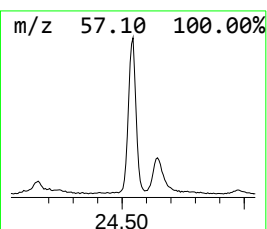
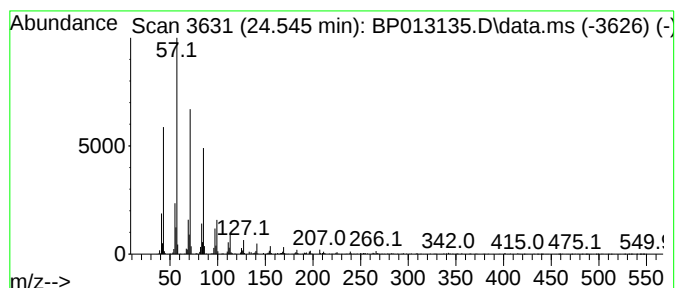
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.545	5.89 ng/ul	1853730	Perylene-d12	23.921

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triacontane	422	C30H62	000638-68-6	94
2			Hexacosane	366	C26H54	000630-01-3	94
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
4			Nonacosane, 2-methyl-	422	C30H62	001560-75-4	93
5			Tetratetracontane	619	C44H90	007098-22-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
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Sample : N6006-13
Misc :
ALS Vial : 14 Sample Multiplier: 1

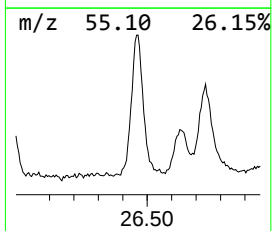
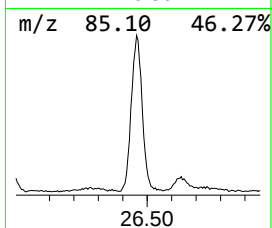
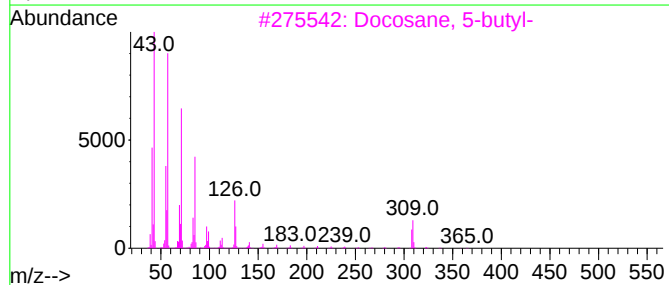
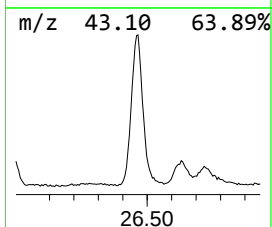
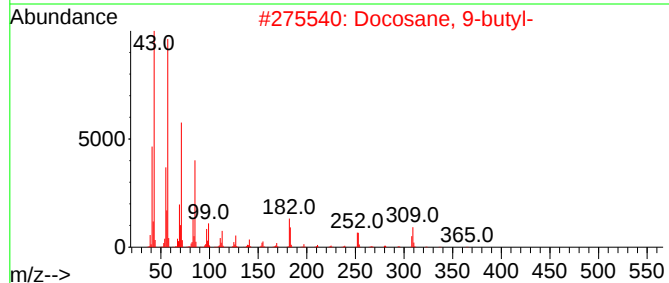
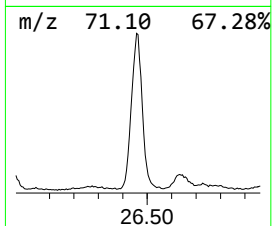
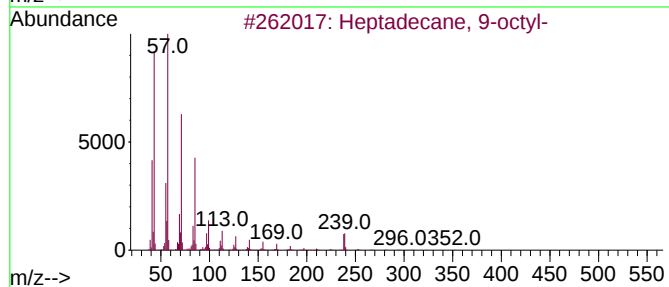
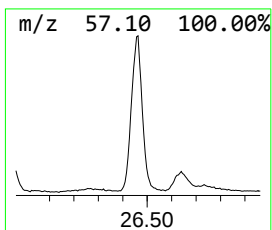
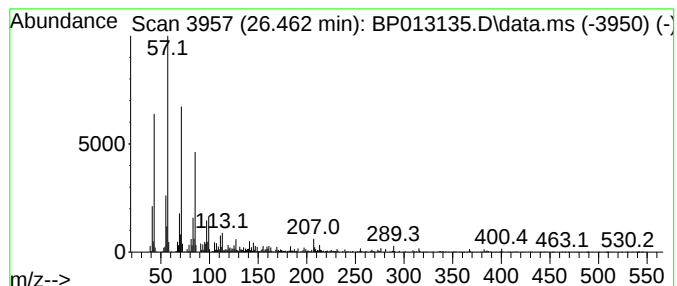
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
26.462	12.33 ng/ul	3882610	Perylene-d12	23.921	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
2	Docosane, 9-butyl-	366	C26H54	055282-14-9	93
3	Docosane, 5-butyl-	366	C26H54	055282-16-1	93
4	2-Bromo dodecane	248	C12H25Br	013187-99-0	91
5	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013135.D
Acq On : 19 Dec 2022 20:04
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Sample : N6006-13
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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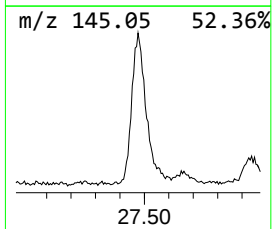
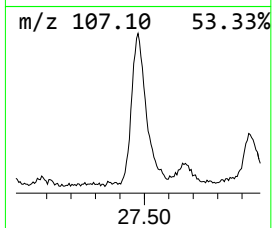
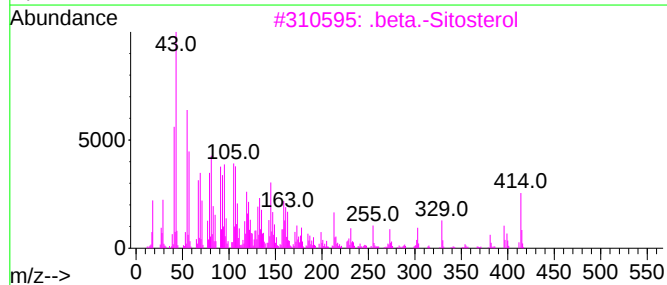
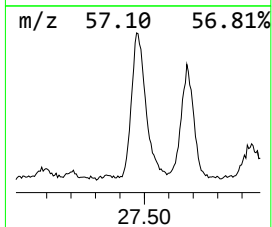
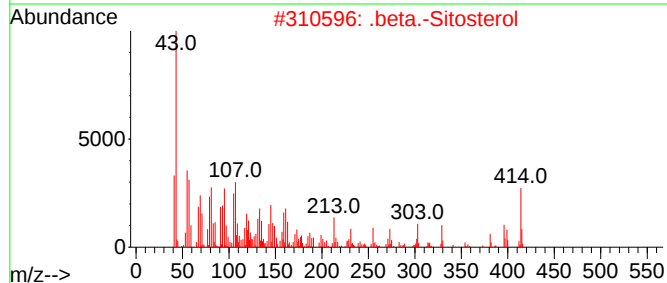
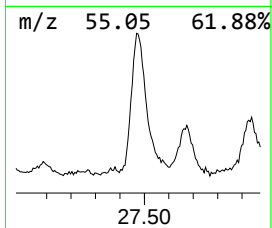
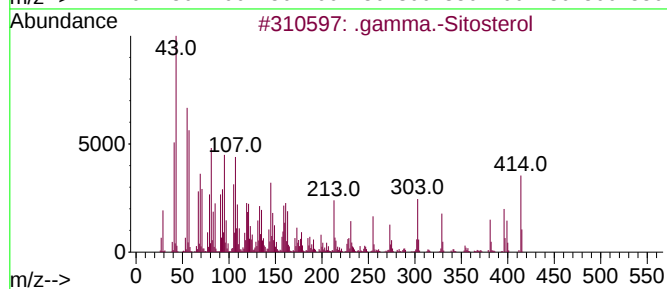
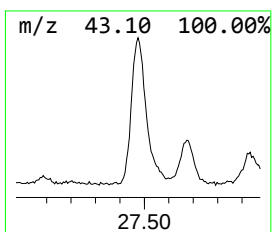
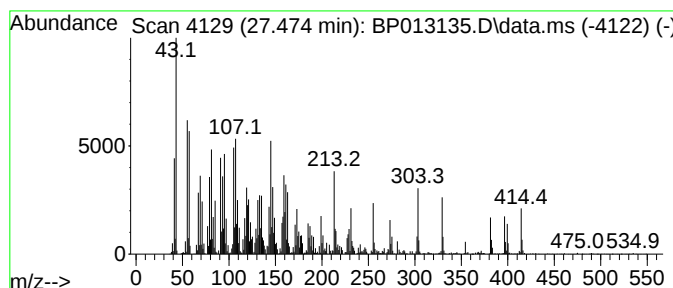
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 .gamma.-Sitosterol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.474	18.60 ng/ul	5855800	Perylene-d12	23.921

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	gamma.-Sitosterol	414	C29H50O	000083-47-6	99	
2	.	beta.-Sitosterol	414	C29H50O	000083-46-5	91	
3	.	beta.-Sitosterol	414	C29H50O	000083-46-5	86	
4	.	gamma.-Sitosterol	414	C29H50O	000083-47-6	53	
5		Campesterol	400	C28H48O	000474-62-4	46	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013135.D
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Misc :
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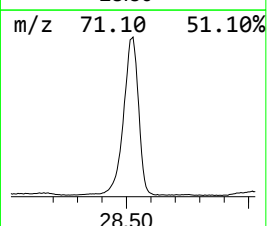
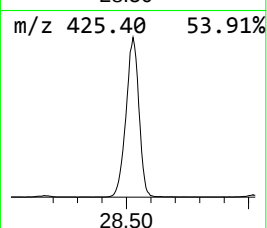
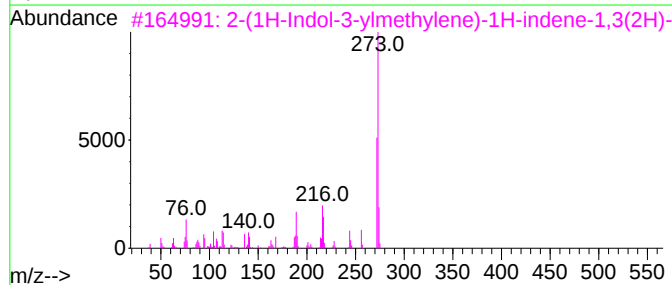
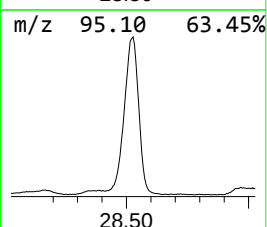
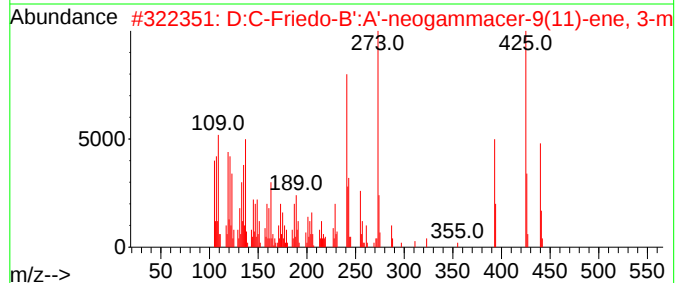
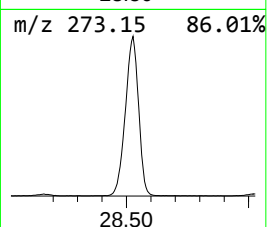
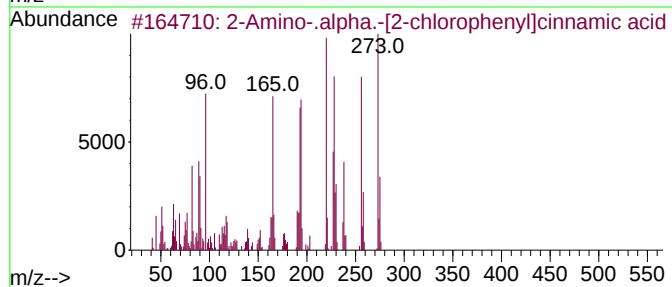
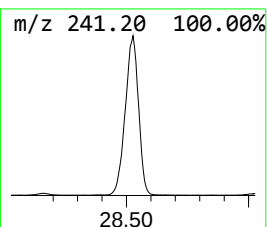
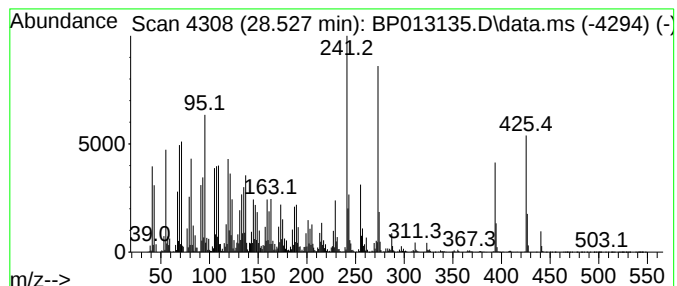
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 2-Amino-.alpha.-[2-chloroph... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
28.527	130.57 ng/ul	41113700	Perylene-d12	23.921		
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Amino-.alpha.-[2-chlorophenyl]...	273	C15H12ClNO2	1000254-08-5	59
2		D:C-Friedo-B':A'-neogammacer-9(1...	440	C31H52O	004555-56-0	59
3		2-(1H-Indol-3-ylmethylene)-1H-in...	273	C18H11NO2	031060-64-7	25
4		1-Phenanthrenecarboxylic acid, 7...	316	C21H32O2	001686-62-0	18
5		2-{[(4-Methyl-2-pyridinyl)amino]...	256	C13H16N4Si	1010494-07-4	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013135.D
Acq On : 19 Dec 2022 20:04
Operator : CG/JU
Sample : N6006-13
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ7

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
2-Pentanone, 4-...	5.040	6.5	ng/ul	1084340	1	7.957	3321210	20.0
Benzophenone	15.957	13.2	ng/ul	4457670	3	14.598	6778880	20.0
n-Hexadecanoic ...	18.227	7.4	ng/ul	2911360	4	17.351	7845930	20.0
Octadecanoic acid	19.457	3.9	ng/ul	1273150	5	21.439	6575450	20.0
3,5-di-tert-But...	20.180	4.1	ng/ul	1337300	5	21.439	6575450	20.0
(DEL) Alkane: C...	21.133	8.1	ng/ul	2671440	5	21.439	6575450	20.0
(DEL) Alkane: S...	22.074	2.9	ng/ul	945660	5	21.439	6575450	20.0
(DEL) Alkane: S...	24.545	5.9	ng/ul	1853730	6	23.921	6297510	20.0
(DEL) Alkane: S...	26.462	12.3	ng/ul	3882610	6	23.921	6297510	20.0
.gamma.-Sitosterol	27.474	18.6	ng/ul	5855800	6	23.921	6297510	20.0
2-Amino-.alpha....	28.527	130.6	ng/ul	41113700	6	23.921	6297510	20.0