

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013140.D
 Acq On : 19 Dec 2022 22:54
 Operator : CG/JU
 Sample : N6006-18
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXQ6

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: BP013140.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	36	rVB	170480	237449	0.55%	0.093%
2	3.617	69	73	86	rBV	775934	1033587	2.40%	0.403%
3	3.770	97	99	113	rBV	661494	1009965	2.34%	0.394%
4	3.999	133	138	147	rVB	72857	116356	0.27%	0.045%
5	4.099	151	155	161	rVB	41963	60762	0.14%	0.024%
6	4.323	189	193	200	rVB	82731	110288	0.26%	0.043%
7	4.740	259	264	270	rVB	37185	57718	0.13%	0.023%
8	5.040	308	315	325	rVB2	184167	302946	0.70%	0.118%
9	7.111	659	667	678	rBV	2988102	4689112	10.88%	1.830%
10	7.281	688	696	714	rVB	2619355	3960562	9.19%	1.546%
11	7.481	723	730	739	rBV	3016662	4773352	11.08%	1.863%
12	7.958	802	811	817	rBV	2074768	3310220	7.68%	1.292%
13	8.664	922	931	942	rBV	3546675	5543855	12.87%	2.164%
14	8.787	946	952	958	rVB	111148	190293	0.44%	0.074%
15	9.122	1002	1009	1018	rBV	2986900	4887604	11.34%	1.907%
16	9.528	1070	1078	1087	rBV	88056	139706	0.32%	0.055%
17	9.846	1124	1132	1147	rBV	2507882	4156202	9.65%	1.622%
18	9.987	1150	1156	1161	rVV4	23215	52316	0.12%	0.020%
19	10.328	1208	1214	1218	rBV	128277	198555	0.46%	0.077%
20	10.387	1218	1224	1243	rVB	3970176	6475094	15.03%	2.527%
21	10.769	1280	1289	1296	rBV	3054615	5058508	11.74%	1.974%
22	10.905	1305	1312	1333	rVV	2708839	4766478	11.06%	1.860%
23	11.552	1416	1422	1428	rBV5	25934	56942	0.13%	0.022%
24	12.175	1524	1528	1534	rVB	78592	110219	0.26%	0.043%
25	12.352	1553	1558	1563	rBV	68800	108576	0.25%	0.042%
26	12.646	1600	1608	1614	rVB2	21594	44391	0.10%	0.017%
27	13.410	1728	1738	1745	rBV	139353	222690	0.52%	0.087%
28	13.487	1745	1751	1754	rBV6	24183	44735	0.10%	0.017%
29	13.563	1761	1764	1771	rVB6	19477	43424	0.10%	0.017%
30	13.916	1820	1824	1830	rBV6	36500	63240	0.15%	0.025%
31	14.004	1831	1839	1847	rVV	7140996	9705761	22.52%	3.788%
32	14.069	1847	1850	1854	rVV	45295	63019	0.15%	0.025%
33	14.116	1854	1858	1862	rVB4	31634	50724	0.12%	0.020%
34	14.293	1881	1888	1903	rBV	8100991	11890563	27.59%	4.640%
35	14.481	1915	1920	1926	rVB	106793	146887	0.34%	0.057%
36	14.599	1929	1940	1947	rBV	4818964	7023748	16.30%	2.741%

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Smoothing : OFF

Filtering: 5

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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	14.657	1947	1950	1965	rVB3	47309	96448	0.22%	0.038%
38	14.793	1966	1973	1990	rBV	2371880	3796280	8.81%	1.482%
39	14.946	1995	1999	2003	rVV5	32605	55411	0.13%	0.022%
40	15.010	2003	2010	2017	rVB5	92820	192084	0.45%	0.075%
41	15.434	2077	2082	2086	rVB2	37610	59928	0.14%	0.023%
42	15.593	2101	2109	2116	rBV	10463597	14907069	34.59%	5.818%
43	15.704	2122	2128	2138	rBV	2994204	4118819	9.56%	1.607%
44	15.828	2143	2149	2157	rVB2	51616	99938	0.23%	0.039%
45	15.957	2165	2171	2181	rVB	2883452	3878058	9.00%	1.513%
46	16.298	2224	2229	2231	rBV3	44253	66259	0.15%	0.026%
47	16.522	2264	2267	2272	rVB2	31750	43760	0.10%	0.017%
48	16.734	2299	2303	2306	rBV4	42118	52376	0.12%	0.020%
49	17.004	2344	2349	2357	rVB5	42271	76555	0.18%	0.030%
50	17.234	2382	2388	2396	rBV3	131180	218294	0.51%	0.085%
51	17.298	2396	2399	2402	rVV3	55152	73104	0.17%	0.029%
52	17.351	2402	2408	2413	rVV	5786829	8408525	19.51%	3.282%
53	17.393	2413	2415	2419	rVV	441230	572459	1.33%	0.223%
54	17.451	2419	2425	2434	rVV2	10537726	14991893	34.79%	5.851%
55	17.540	2437	2440	2447	rVB6	76969	125947	0.29%	0.049%
56	17.751	2474	2476	2481	rVB	49469	60632	0.14%	0.024%
57	17.840	2487	2491	2494	rBV2	45826	64387	0.15%	0.025%
58	18.010	2516	2520	2523	rBV	71102	76339	0.18%	0.030%
59	18.110	2532	2537	2542	rBV2	107194	170508	0.40%	0.067%
60	18.163	2542	2546	2551	rBV	271735	386524	0.90%	0.151%
61	18.228	2551	2557	2567	rBV	2480319	3536441	8.21%	1.380%
62	18.510	2601	2605	2609	rBV4	89513	136513	0.32%	0.053%
63	18.551	2610	2612	2618	rVB3	51911	69017	0.16%	0.027%
64	18.634	2623	2626	2631	rBV3	67455	92962	0.22%	0.036%
65	18.698	2634	2637	2641	rVB2	70329	89869	0.21%	0.035%
66	18.863	2662	2665	2668	rBV3	51126	64585	0.15%	0.025%
67	18.928	2674	2676	2682	rVB6	63725	78647	0.18%	0.031%
68	19.104	2701	2706	2711	rBV4	261032	476852	1.11%	0.186%
69	19.257	2729	2732	2738	rVB2	186317	267820	0.62%	0.105%
70	19.322	2738	2743	2744	rBV3	286187	415920	0.97%	0.162%
71	19.357	2745	2749	2757	rVV	1105957	1654052	3.84%	0.646%
72	19.451	2761	2765	2778	rVV	624322	950219	2.21%	0.371%
73	19.681	2798	2804	2816	rBV	12801619	18419216	42.74%	7.188%
74	20.192	2888	2891	2898	rVB2	204096	264539	0.61%	0.103%
75	21.128	3046	3050	3055	rBV	2294081	3312940	7.69%	1.293%
76	21.433	3097	3102	3107	rBV	5024676	7234271	16.79%	2.823%

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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

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Title : SVOA CALIBRATION

77	23.133	3386	3391	3399	rBV2	1386527	2813020	6.53%	1.098%
78	23.757	3490	3497	3513	rVB2	6209585	13241273	30.73%	5.168%
79	23.916	3518	3524	3543	rVB2	2840443	6284004	14.58%	2.452%
80	24.539	3624	3630	3639	rVB	1603364	3381707	7.85%	1.320%
81	25.392	3768	3775	3788	rVB5	1552184	4634729	10.76%	1.809%
82	26.451	3948	3955	3977	rVB3	1662984	6084615	14.12%	2.375%
83	27.463	4120	4127	4146	rVB4	1702708	6343916	14.72%	2.476%
84	28.515	4293	4306	4326	rVB2	10756194	43091668	100.00%	16.817%

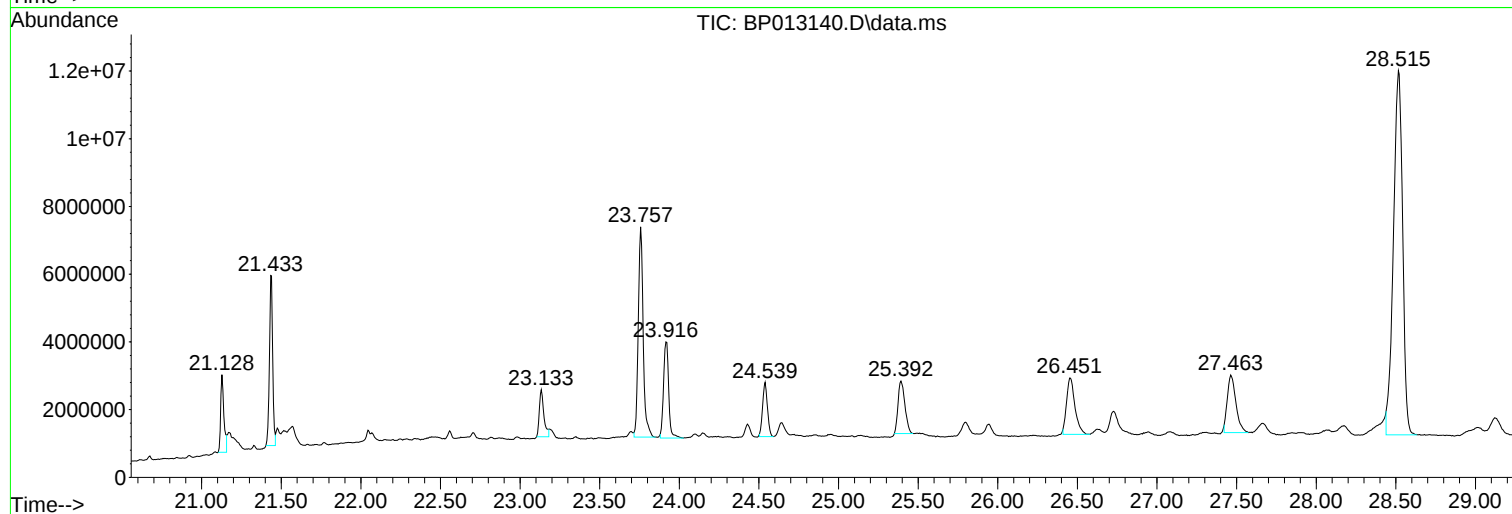
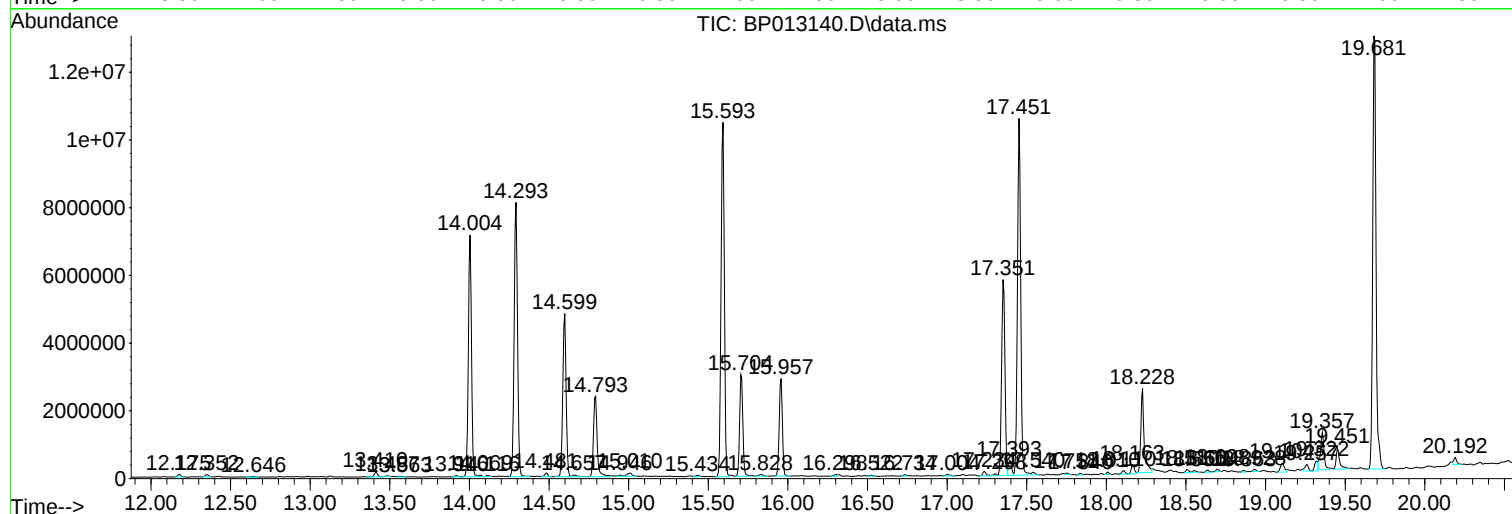
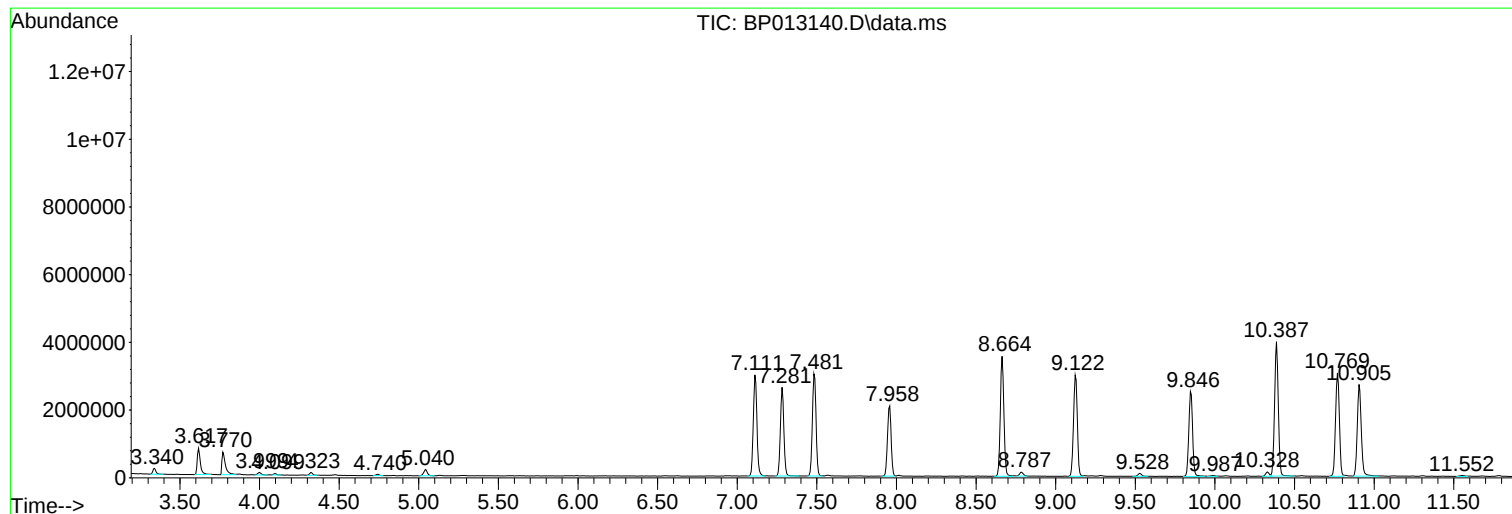
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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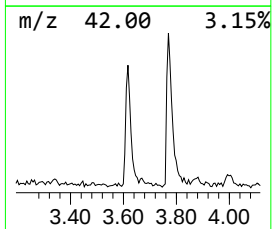
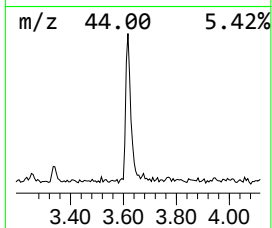
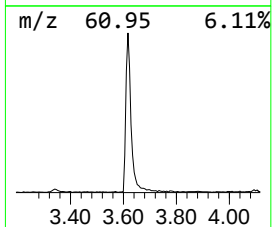
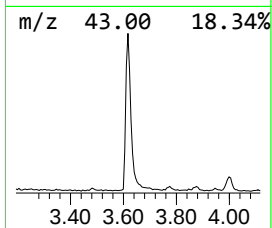
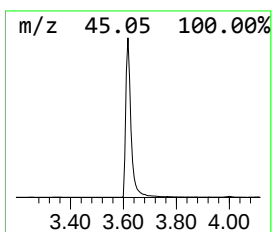
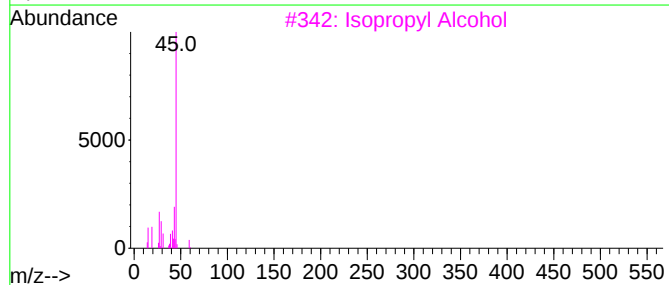
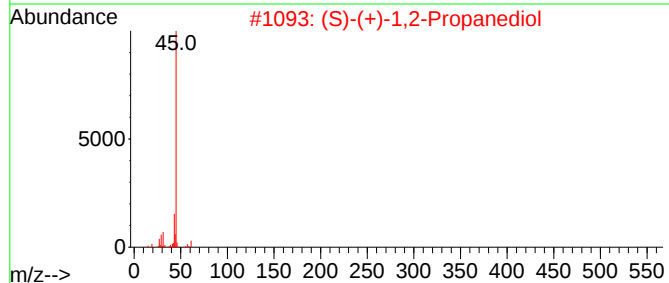
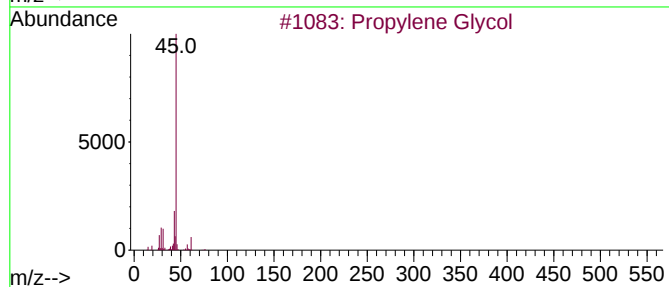
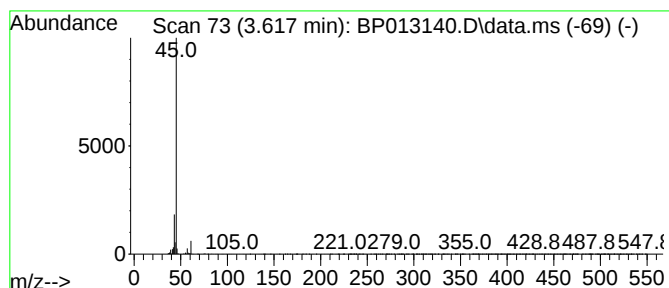
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.617	6.24 ng/ul	1033590	1,4-Dichlorobenzene-d4	7.958

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	90
2			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	64
3			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
4			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
5			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40



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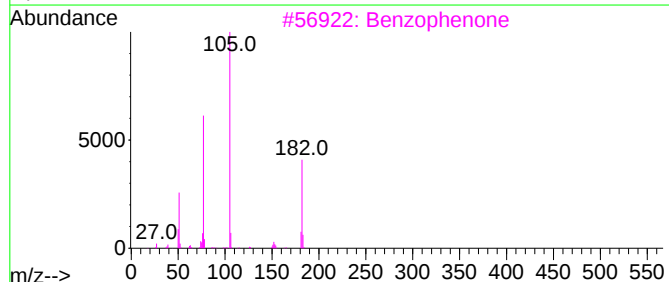
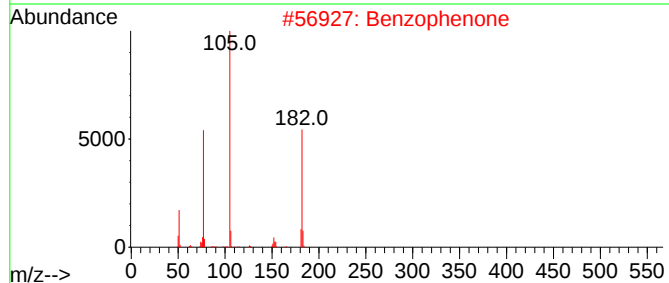
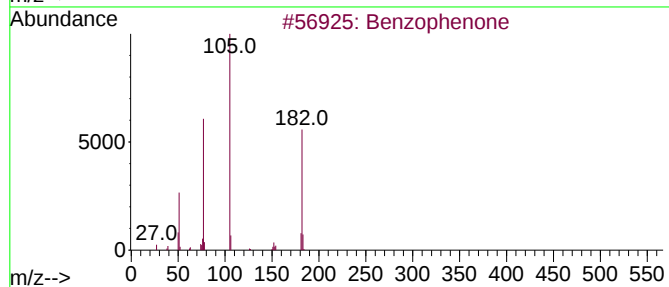
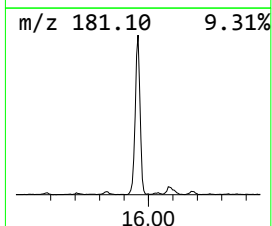
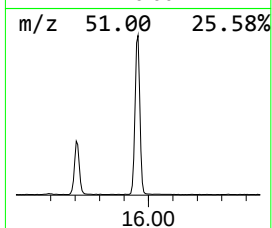
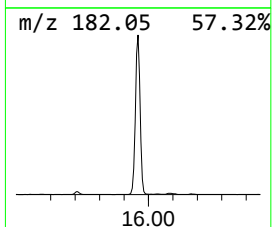
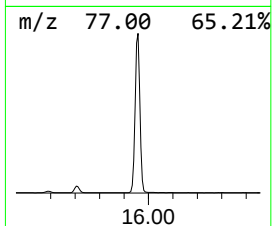
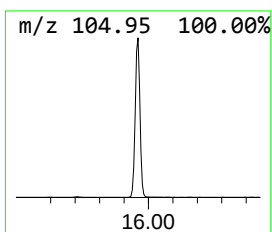
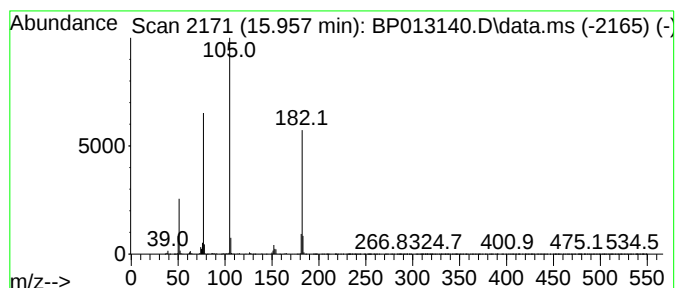
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.957	11.04 ng/ul	3878060	Acenaphthene-d10	14.599

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	93
5			Benzophenone	182	C13H10O	000119-61-9	91



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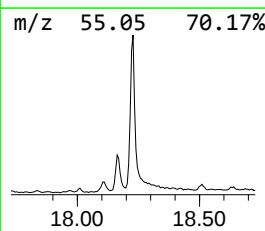
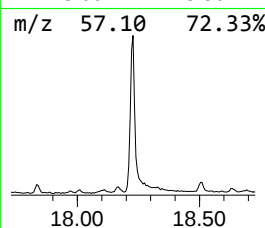
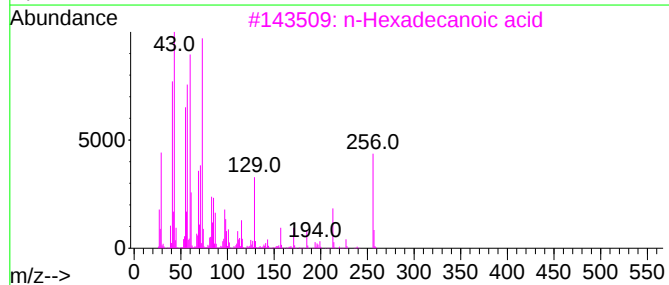
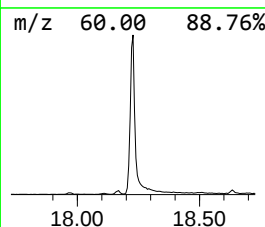
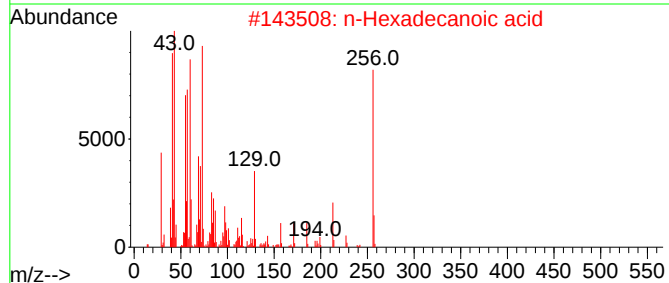
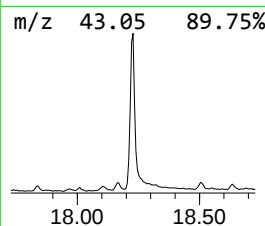
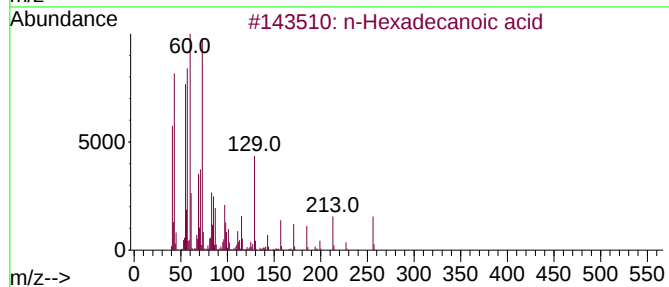
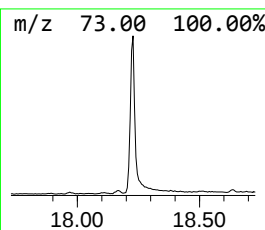
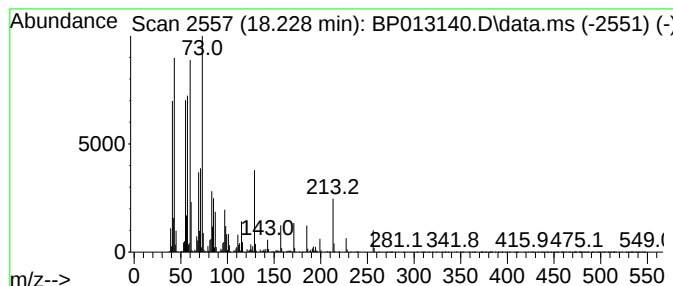
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.228	8.41 ng/ul	3536440	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	98
3	n-Hexadecanoic acid			256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid			256	C16H32O2	000057-10-3	97
5	n-Hexadecanoic acid			256	C16H32O2	000057-10-3	95



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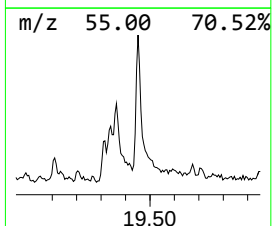
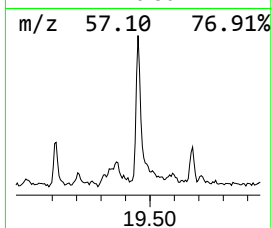
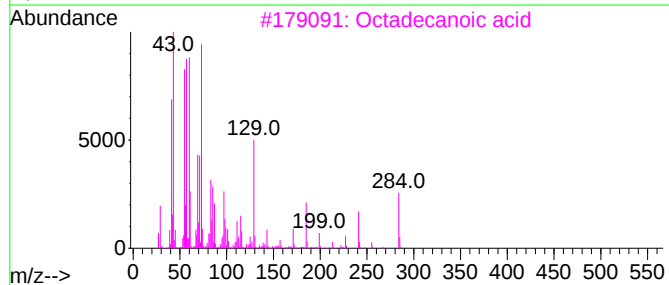
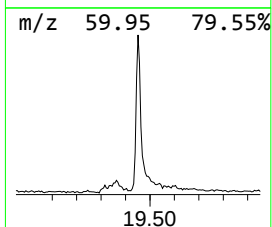
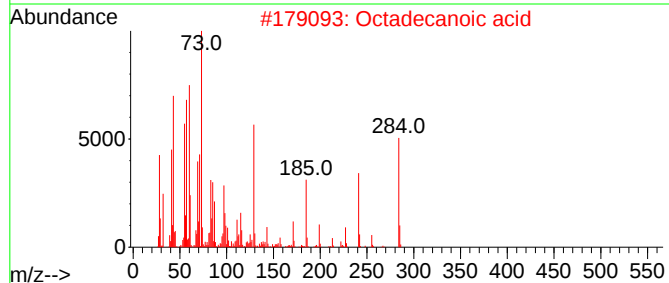
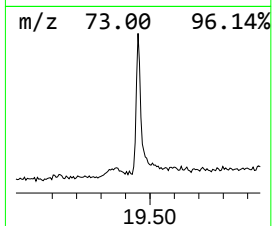
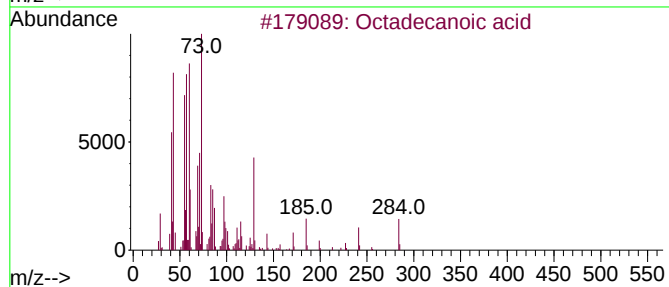
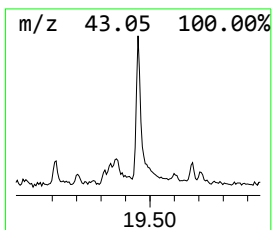
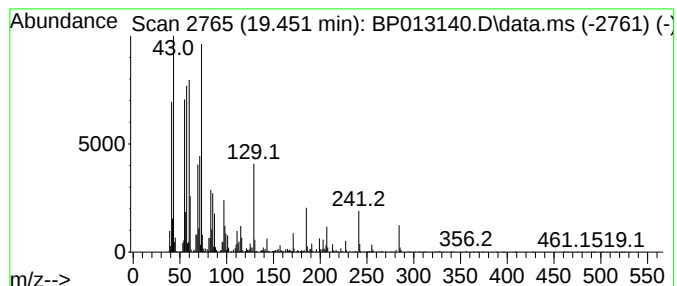
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.451	2.63 ng/ul	950219	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	94
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013140.D
Acq On : 19 Dec 2022 22:54
Operator : CG/JU
Sample : N6006-18
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ6

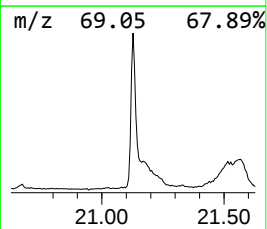
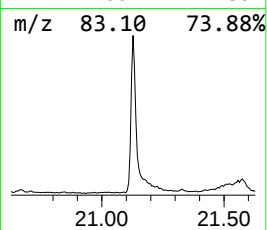
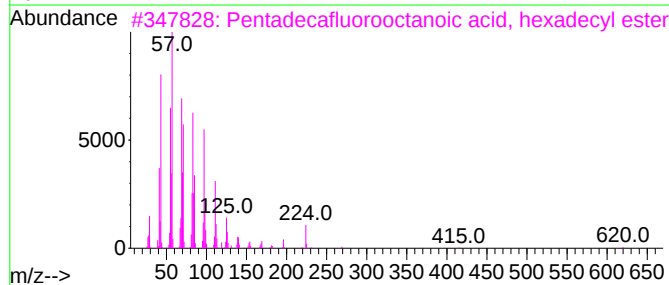
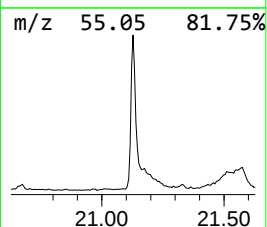
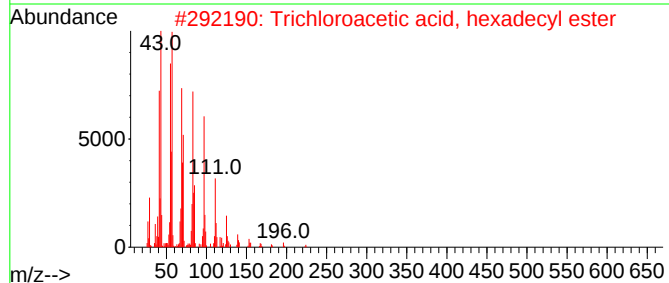
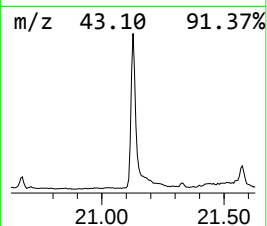
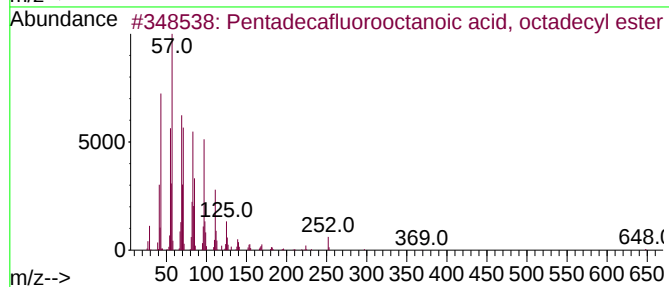
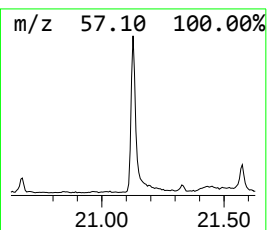
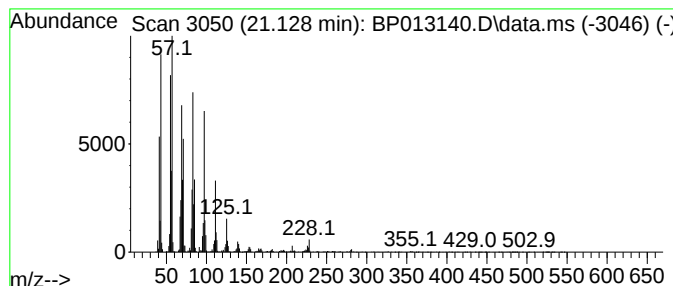
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Pentadecafluorooctanoic aci... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.128	9.16 ng/ul	3312940	Chrysene-d12	21.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
2			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3			Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	94
4			Cyclopentadecane	210	C15H30	000295-48-7	93
5			1-Heneicosanol	312	C21H44O	015594-90-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013140.D
Acq On : 19 Dec 2022 22:54
Operator : CG/JU
Sample : N6006-18
Misc :
ALS Vial : 19 Sample Multiplier: 1

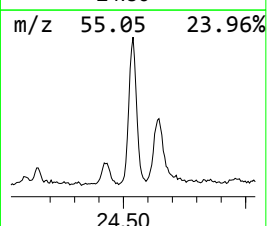
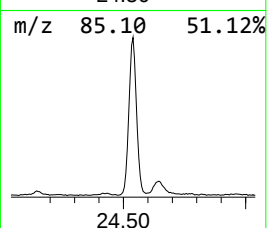
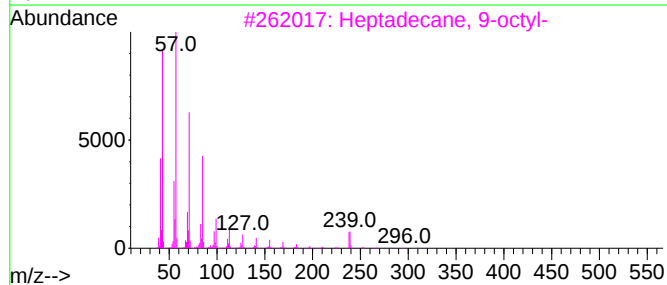
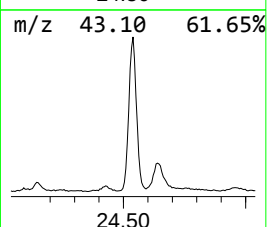
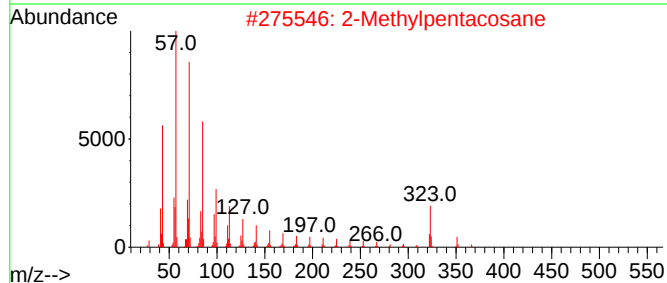
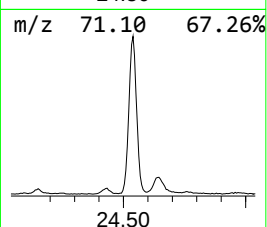
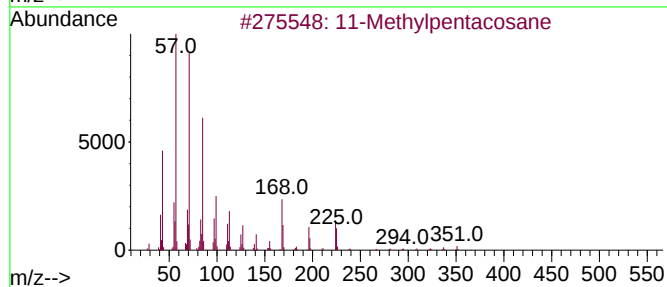
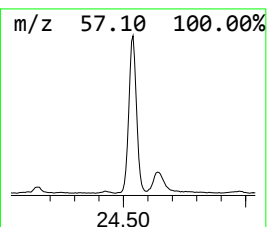
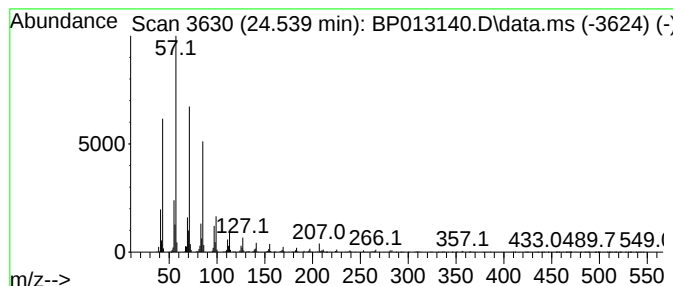
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.539	10.76 ng/ul	3381710	Perylene-d12	23.916	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11-Methylpentacosane	366	C26H54	015689-71-1	93
2	2-Methylpentacosane	366	C26H54	000629-87-8	93
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013140.D
Acq On : 19 Dec 2022 22:54
Operator : CG/JU
Sample : N6006-18
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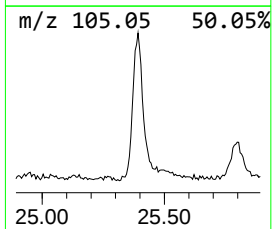
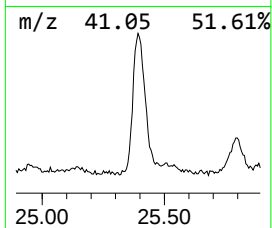
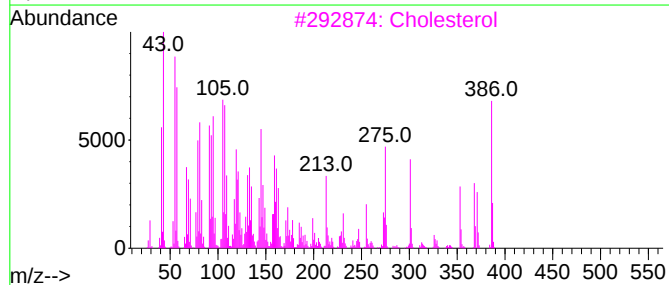
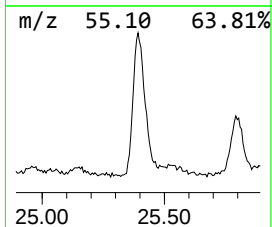
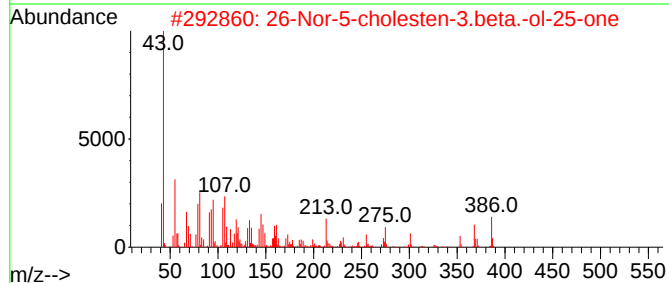
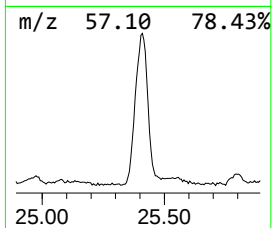
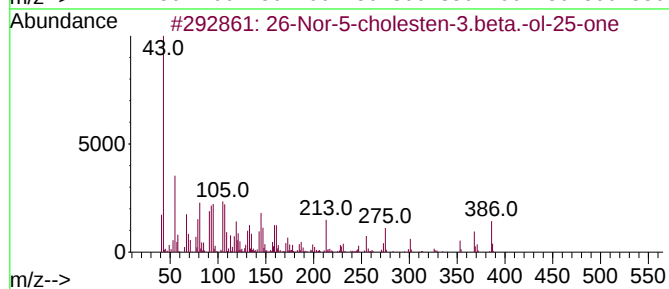
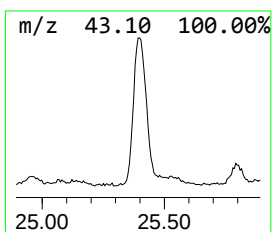
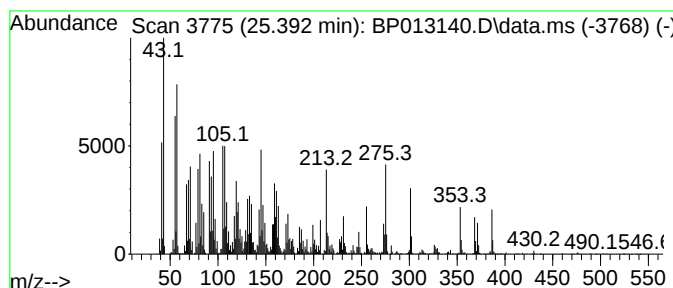
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 26-Nor-5-cholesten-3.beta.-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
25.392	14.75 ng/ul	4634730	Perylene-d12	23.916	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	99
2	26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	99
3	Cholesterol	386	C27H46O	000057-88-5	99
4	26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	78
5	Lathosterol	386	C27H46O	000080-99-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013140.D
Acq On : 19 Dec 2022 22:54
Operator : CG/JU
Sample : N6006-18
Misc :
ALS Vial : 19 Sample Multiplier: 1

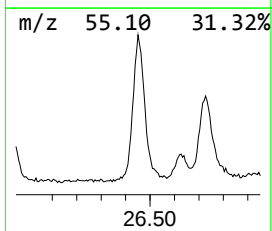
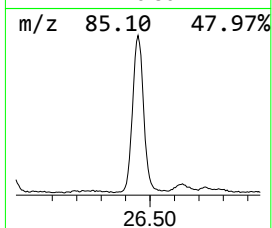
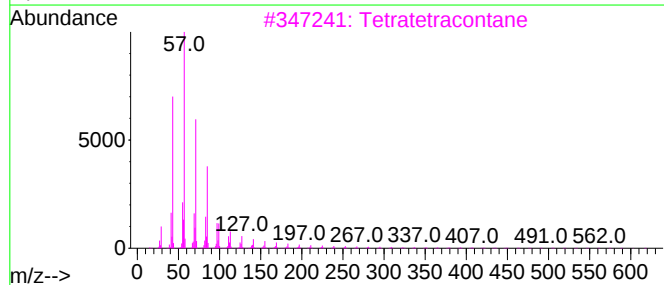
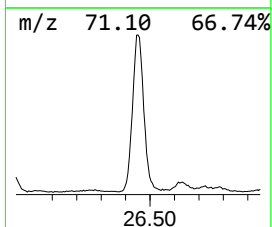
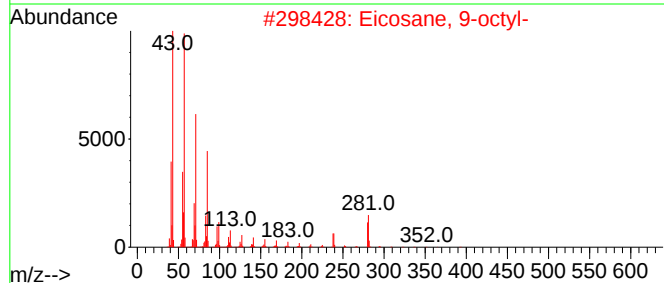
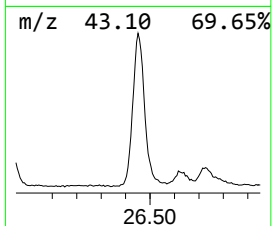
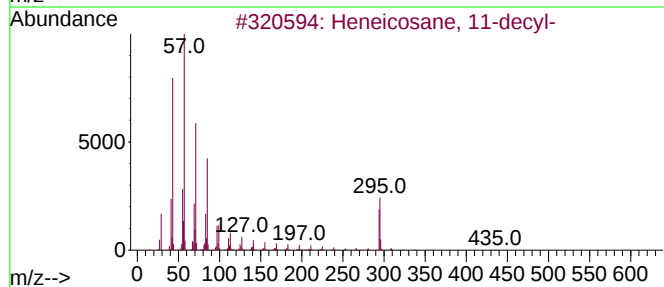
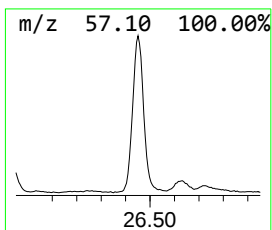
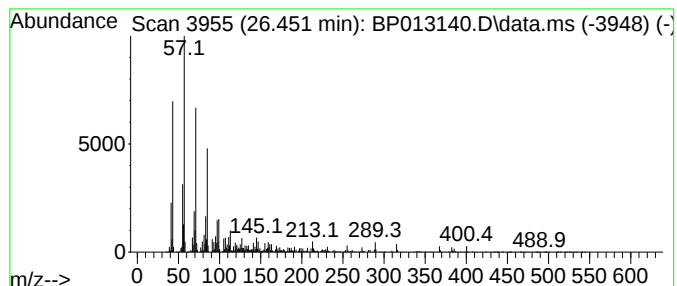
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
26.451	19.37 ng/ul	6084620	Perylene-d12	23.916	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	98
2	Eicosane, 9-octyl-	394	C28H58	013475-77-9	93
3	Tetratetracontane	619	C44H90	007098-22-8	93
4	Tetracosane, 11-decyl-	479	C34H70	055429-84-0	90
5	Tricosane	324	C23H48	000638-67-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013140.D
Acq On : 19 Dec 2022 22:54
Operator : CG/JU
Sample : N6006-18
Misc :
ALS Vial : 19 Sample Multiplier: 1

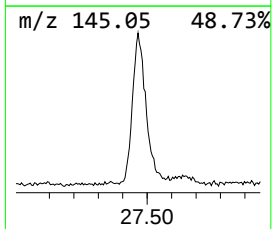
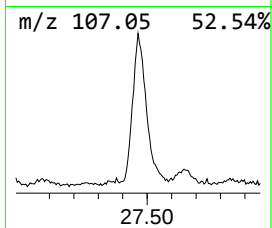
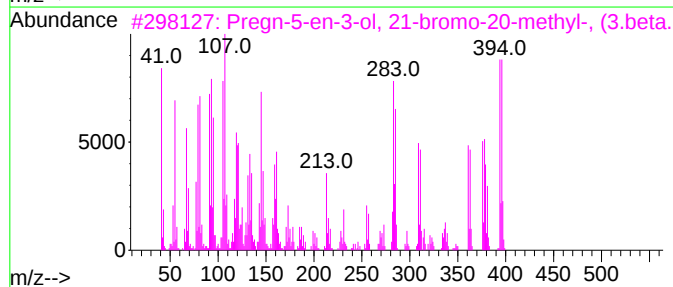
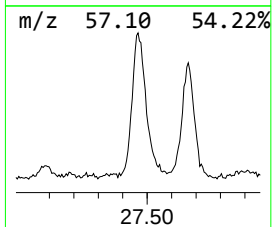
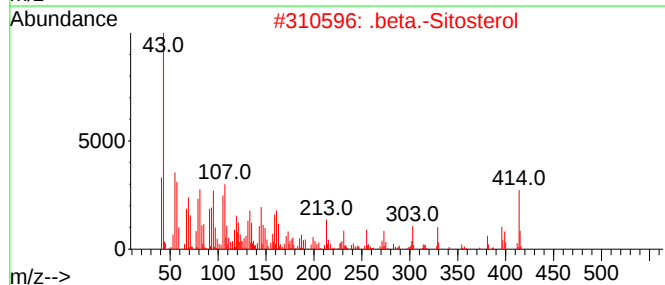
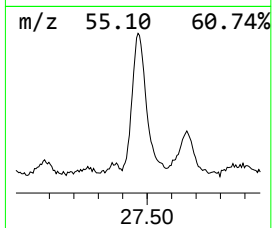
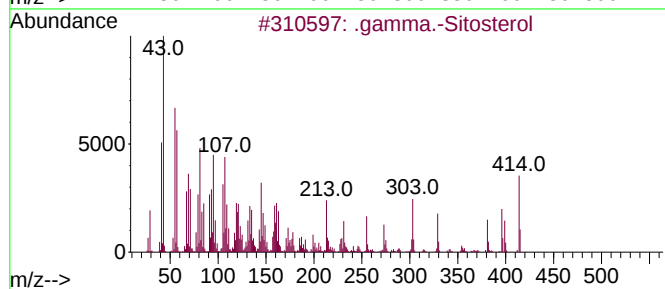
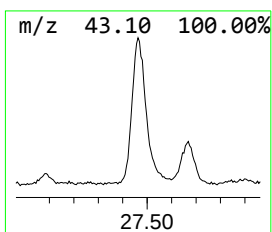
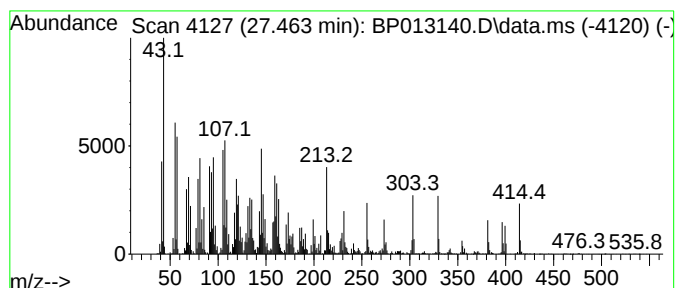
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 .gamma.-Sitosterol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
27.463	20.19 ng/ul	6343920	Perylene-d12	23.916	
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	94
3	Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	91
4	.beta.-Sitosterol	414	C29H50O	000083-46-5	87
5	.gamma.-Sitosterol	414	C29H50O	000083-47-6	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013140.D
Acq On : 19 Dec 2022 22:54
Operator : CG/JU
Sample : N6006-18
Misc :
ALS Vial : 19 Sample Multiplier: 1

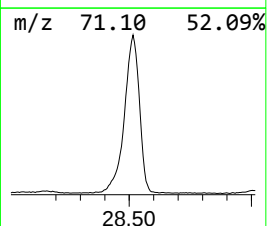
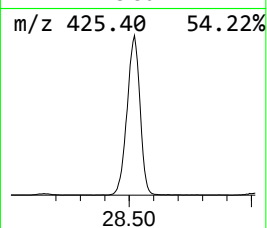
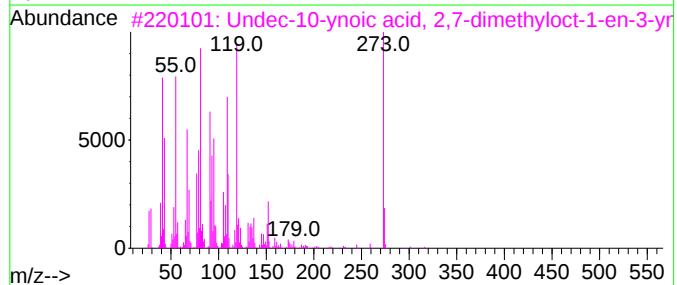
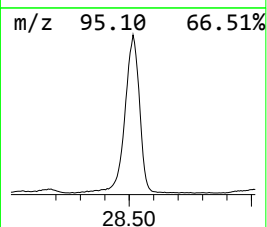
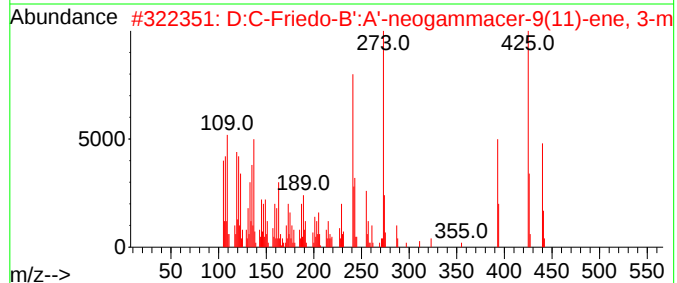
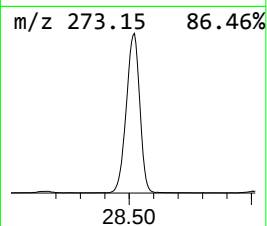
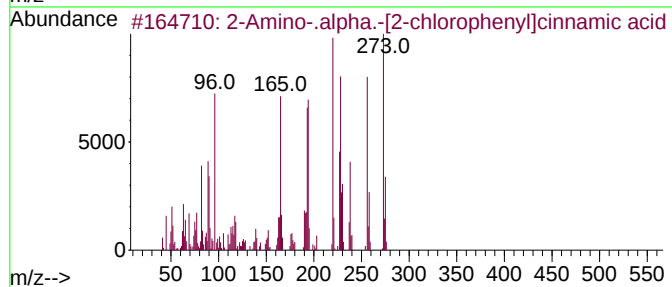
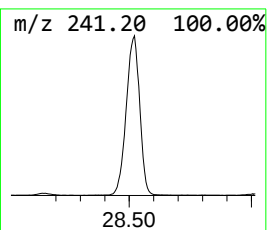
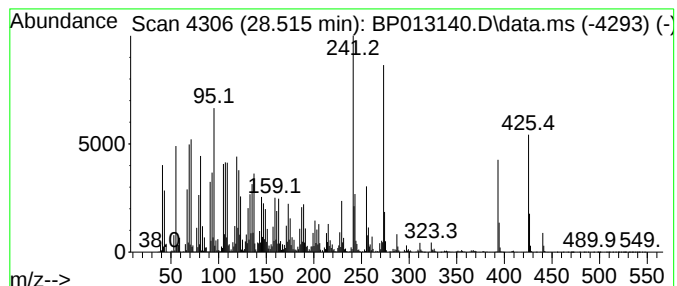
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 2-Amino-.alpha.-[2-chloroph... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
28.515	137.15 ng/ul	43091700	Perylene-d12			23.916
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	43
3		Undec-10-ynoic acid, 2,7-dimethy...	316	C21H32O2	1000406-96-7	22
4		Pyrazolo[1,5-a]pyrimidine, 6,7-d...	273	C18H15N3	186956-30-9	15
5		2-[(4-Methyl-2-pyridinyl)amino]...	256	C13H16N4Si	1010494-07-4	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013140.D
Acq On : 19 Dec 2022 22:54
Operator : CG/JU
Sample : N6006-18
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXQ6

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
Propylene Glycol	3.617	6.2	ng/ul	1033590	1	7.958	3310220	20.0
Benzophenone	15.957	11.0	ng/ul	3878060	3	14.599	7023750	20.0
n-Hexadecanoic ...	18.228	8.4	ng/ul	3536440	4	17.351	8408530	20.0
Octadecanoic acid	19.451	2.6	ng/ul	950219	5	21.439	7234270	20.0
Pentadecafluoro...	21.128	9.2	ng/ul	3312940	5	21.439	7234270	20.0
(DEL) Alkane: S...	24.539	10.8	ng/ul	3381710	6	23.916	6284000	20.0
26-Nor-5-choles...	25.392	14.8	ng/ul	4634730	6	23.916	6284000	20.0
(DEL) Alkane: S...	26.451	19.4	ng/ul	6084620	6	23.916	6284000	20.0
.gamma.-Sitosterol	27.463	20.2	ng/ul	6343920	6	23.916	6284000	20.0
2-Amino-.alpha....	28.515	137.2	ng/ul	43091700	6	23.916	6284000	20.0