

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
 Data File : BG059105.D
 Acq On : 20 Sep 2023 12:01
 Operator : MA/JU
 Sample : 04378-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BBJG5

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_G\Methods\SFAM-EPA-BG091923.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG059105.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.611	34	38	44	rVB4	10505	17263	0.61%	0.113%
2	4.046	107	112	121	rBV	32691	57264	2.02%	0.376%
3	4.258	144	148	151	rVB2	8877	10845	0.38%	0.071%
4	4.916	254	260	271	rVB2	43659	72150	2.54%	0.473%
5	7.201	644	649	654	rVB2	4458	7690	0.27%	0.050%
6	7.407	678	684	693	rVB2	36205	63397	2.24%	0.416%
7	7.618	713	720	726	rVB	117618	188196	6.64%	1.234%
8	7.812	745	753	762	rVB2	109131	182878	6.45%	1.200%
9	8.188	812	817	822	rVB3	4367	6688	0.24%	0.044%
10	8.300	829	836	843	rBV	109850	184937	6.52%	1.213%
11	8.447	857	861	866	rBV2	3729	6670	0.24%	0.044%
12	8.764	911	915	919	rBV2	4361	6496	0.23%	0.043%
13	8.893	933	937	944	rBV3	3415	5962	0.21%	0.039%
14	8.976	944	951	960	rVV2	105353	181216	6.39%	1.189%
15	9.105	966	973	977	rBV3	2140	3727	0.13%	0.024%
16	9.493	1030	1039	1046	rVB	143211	251075	8.86%	1.647%
17	10.215	1154	1162	1170	rVB	110957	203208	7.17%	1.333%
18	10.427	1194	1198	1203	rBV2	3081	4319	0.15%	0.028%
19	10.503	1207	1211	1216	rVB2	3045	4483	0.16%	0.029%
20	10.738	1242	1251	1262	rBV	175417	316582	11.17%	2.077%
21	11.144	1313	1320	1327	rBV	174146	310386	10.95%	2.036%
22	11.285	1336	1344	1353	rBV	152052	278998	9.84%	1.830%
23	12.706	1581	1586	1588	rVB	2867	3667	0.13%	0.024%
24	13.676	1746	1751	1756	rVB4	7833	11100	0.39%	0.073%
25	13.752	1760	1764	1771	rVB3	4940	7074	0.25%	0.046%
26	14.322	1851	1861	1867	rBV	376577	549068	19.37%	3.602%
27	14.628	1906	1913	1921	rVB	416681	674271	23.78%	4.423%
28	14.810	1940	1944	1948	rBV	3024	3735	0.13%	0.024%
29	14.927	1957	1964	1972	rBV	284965	435462	15.36%	2.856%
30	15.092	1985	1992	2003	rBV2	44263	71787	2.53%	0.471%
31	15.691	2090	2094	2097	rBV	9833	11574	0.41%	0.076%
32	15.914	2125	2132	2138	rBV	561539	853194	30.09%	5.597%
33	15.961	2138	2140	2144	rVB	2471	3125	0.11%	0.020%
34	16.020	2144	2150	2157	rBV	192396	288467	10.18%	1.892%
35	16.555	2235	2241	2243	rBV2	4720	5564	0.20%	0.036%
36	16.608	2248	2250	2257	rBV3	3826	6125	0.22%	0.040%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
 Data File : BG059105.D
 Acq On : 20 Sep 2023 12:01
 Operator : MA/JU
 Sample : 04378-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BBJG5

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_G\Methods\SFAM-EPA-BG091923.MA.M
 Title : SVOA CALIBRATION

37	17.219	2351	2354	2362	rBV3	2661	4809	0.17%	0.032%
38	17.342	2372	2375	2380	rVV3	6117	7146	0.25%	0.047%
39	17.383	2380	2382	2387	rVB2	2224	2861	0.10%	0.019%
40	17.677	2426	2432	2440	rVB	353523	532863	18.80%	3.495%
41	17.777	2442	2449	2458	rBV2	662818	988134	34.85%	6.482%
42	18.088	2499	2502	2506	rVB3	4571	5177	0.18%	0.034%
43	18.159	2510	2514	2517	rBV5	3410	4368	0.15%	0.029%
44	18.494	2567	2571	2579	rBV2	44996	66185	2.33%	0.434%
45	19.111	2674	2676	2678	rBV2	3329	3055	0.11%	0.020%
46	19.234	2695	2697	2698	rBV2	3642	2884	0.10%	0.019%
47	19.393	2722	2724	2727	rBV4	2665	2981	0.11%	0.020%
48	19.622	2756	2763	2765	rBV6	12022	20158	0.71%	0.132%
49	19.687	2765	2774	2776	rVV4	8368	18017	0.64%	0.118%
50	19.757	2782	2786	2793	rBV4	32144	47175	1.66%	0.309%
51	19.898	2806	2810	2814	rBV4	6960	9463	0.33%	0.062%
52	19.969	2818	2822	2827	rVB6	5216	8193	0.29%	0.054%
53	20.051	2830	2836	2842	rVB	767278	1102239	38.88%	7.230%
54	20.503	2909	2913	2915	rBV4	8143	10055	0.35%	0.066%
55	20.580	2924	2926	2927	rBV2	4987	3319	0.12%	0.022%
56	20.674	2939	2942	2945	rBV4	4310	6485	0.23%	0.043%
57	21.003	2996	2998	3002	rVB3	10351	11079	0.39%	0.073%
58	21.537	3086	3089	3098	rBV4	17808	27547	0.97%	0.181%
59	21.996	3161	3167	3174	rVB2	283660	498249	17.57%	3.268%
60	22.119	3184	3188	3194	rVB	22215	29158	1.03%	0.191%
61	22.436	3240	3242	3246	rVB5	7517	8176	0.29%	0.054%
62	22.771	3295	3299	3307	rVB5	20874	41998	1.48%	0.275%
63	23.529	3423	3428	3440	rVB2	28897	62102	2.19%	0.407%
64	23.741	3458	3464	3472	rVB2	44640	92718	3.27%	0.608%
65	23.952	3499	3500	3502	rBV2	4466	3475	0.12%	0.023%
66	24.410	3572	3578	3585	rBV3	30866	66310	2.34%	0.435%
67	24.628	3612	3615	3617	rBV4	5912	6754	0.24%	0.044%
68	25.298	3717	3729	3742	rBV2	366783	1113688	39.28%	7.305%
69	25.462	3751	3757	3762	rBV2	24533	60652	2.14%	0.398%
70	25.544	3763	3771	3783	rVB2	180912	559940	19.75%	3.673%
71	26.719	3965	3971	3972	rBV4	19188	32557	1.15%	0.214%
72	27.419	4075	4090	4109	rBV4	728757	2835045	100.00%	18.597%
73	28.006	4178	4190	4203	rBV6	199982	740267	26.11%	4.856%
74	29.081	4359	4373	4375	rBV8	36350	104382	3.68%	0.685%
75	29.540	4442	4451	4452	rBV7	10338	25047	0.88%	0.164%
76	29.751	4484	4487	4490	rBV4	4706	7362	0.26%	0.048%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059105.D
Acq On : 20 Sep 2023 12:01
Operator : MA/JU
Sample : 04378-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BBJG5

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_G\Methods\SFAM-EPA-BG091923.MA.M
Title : SVOA CALIBRATION

77	30.680	4630	4645	4667	rBV8	149467	781154	27.55%	5.124%
78	31.678	4813	4815	4816	rBV2	4873	3144	0.11%	0.021%

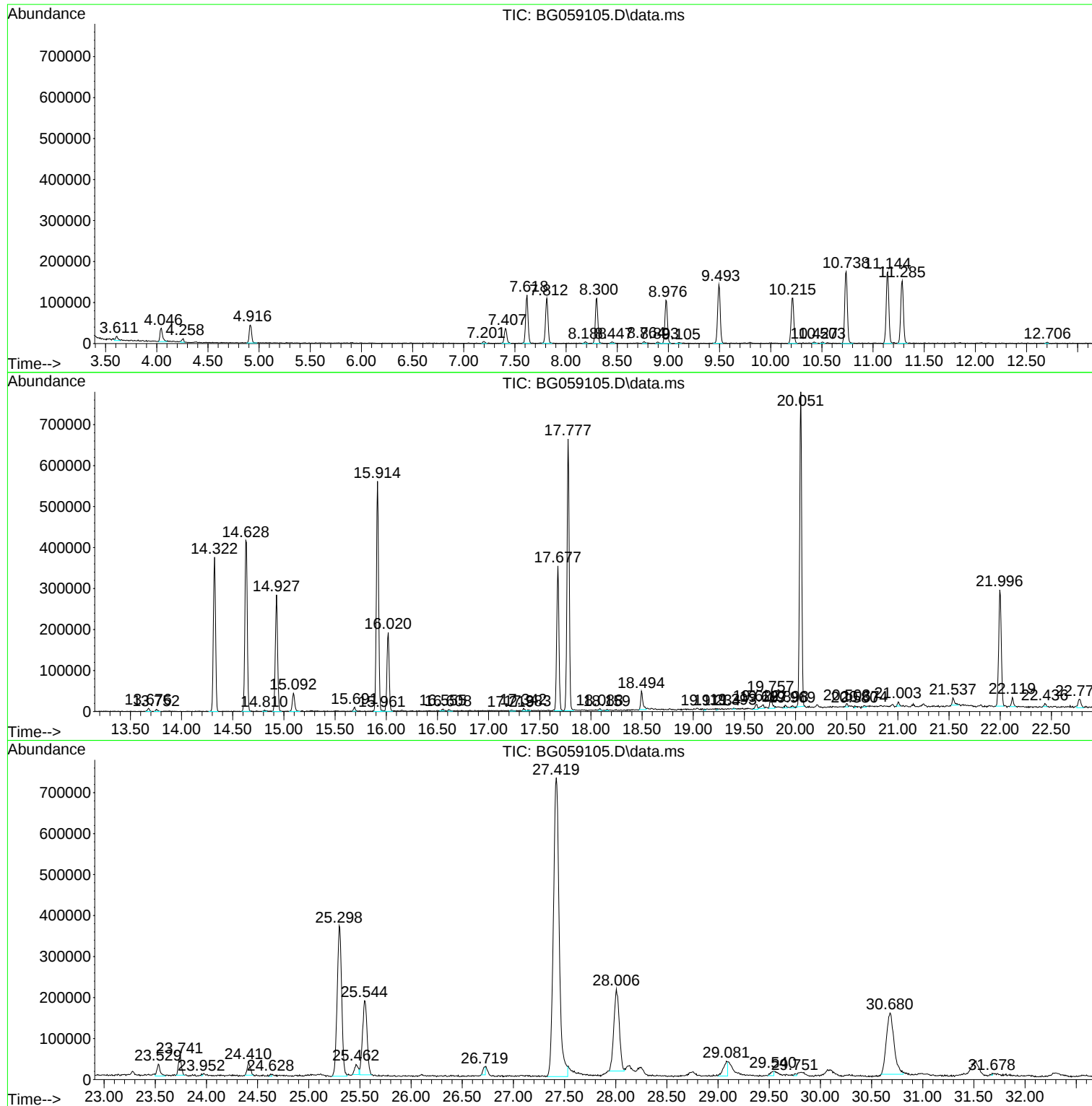
Sum of corrected areas: 15245014

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059105.D
Acq On : 20 Sep 2023 12:01
Operator : MA/JU
Sample : 04378-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BBJG5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059105.D
Acq On : 20 Sep 2023 12:01
Operator : MA/JU
Sample : 04378-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BBJG5

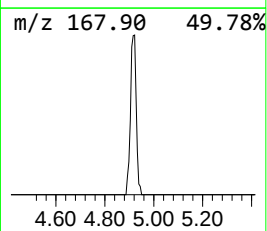
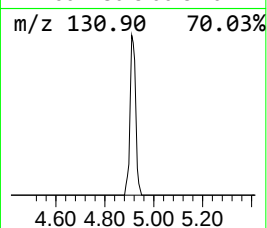
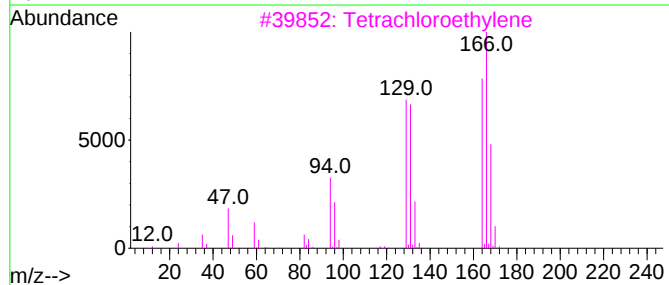
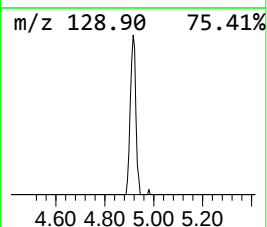
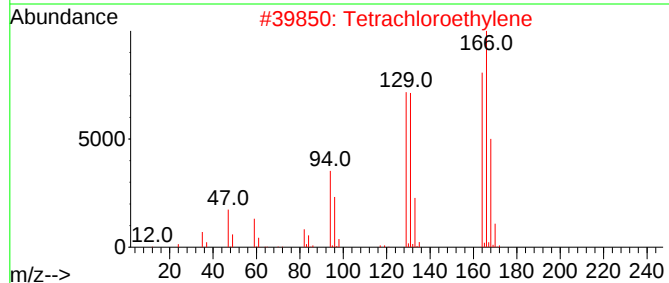
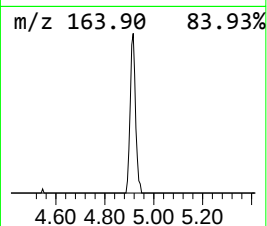
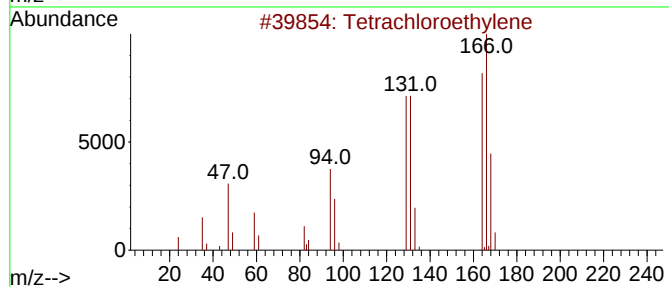
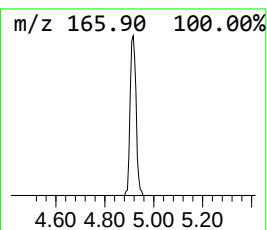
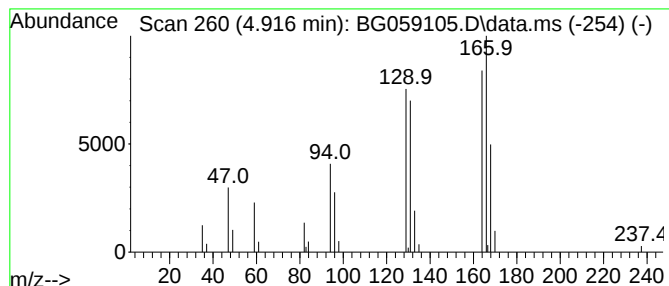
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Tetrachloroethylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.916	7.80 ng/ul	72150	1,4-Dichlorobenzene-d4	8.300

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	97
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	93
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059105.D
Acq On : 20 Sep 2023 12:01
Operator : MA/JU
Sample : 04378-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BBJG5

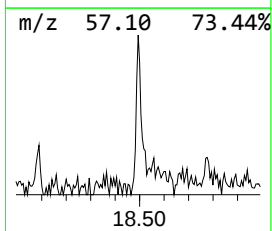
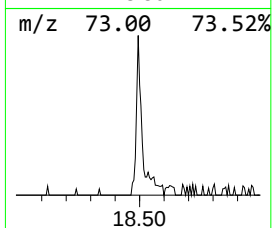
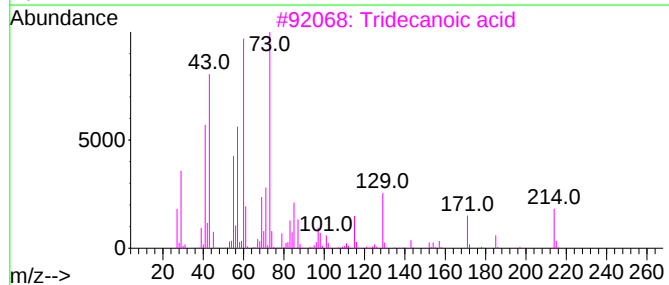
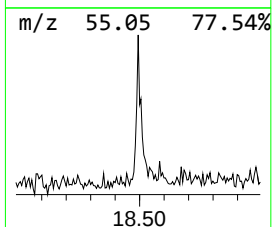
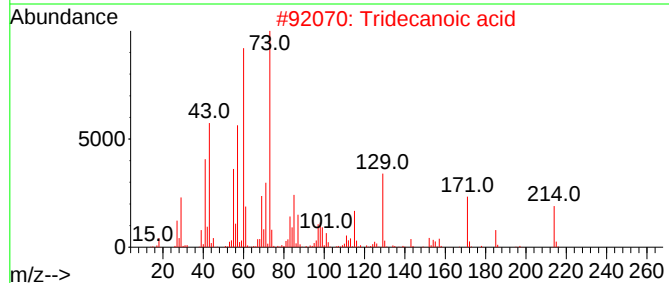
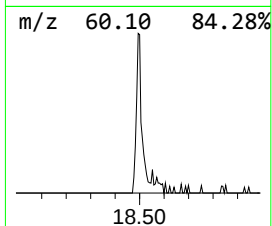
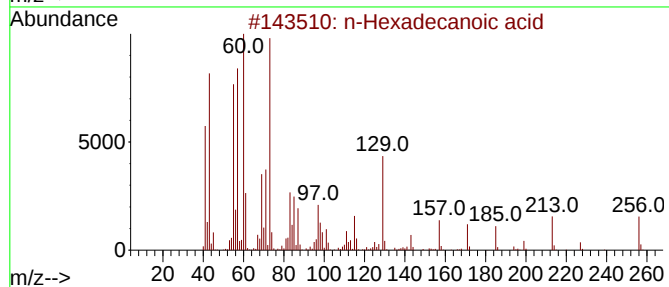
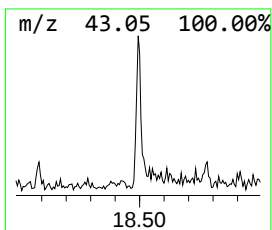
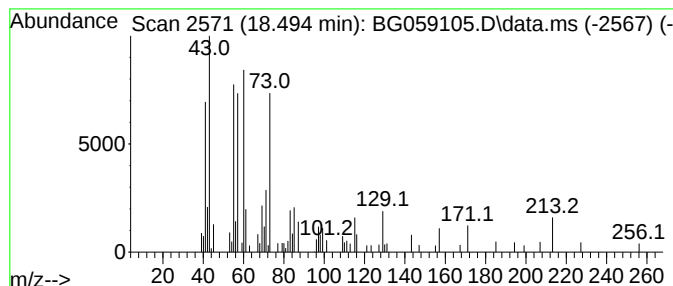
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.494	2.48 ng/ul	66185	Phenanthrene-d10	17.677

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
2			Tridecanoic acid	214	C13H26O2	000638-53-9	80
3			Tridecanoic acid	214	C13H26O2	000638-53-9	74
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	74
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059105.D
Acq On : 20 Sep 2023 12:01
Operator : MA/JU
Sample : 04378-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BBJG5

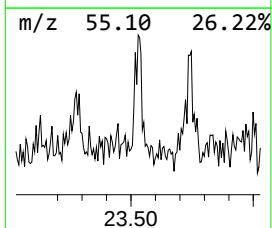
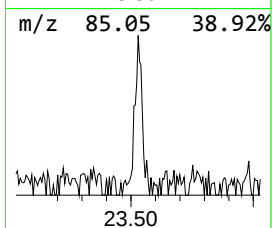
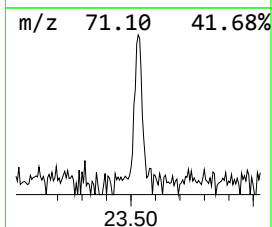
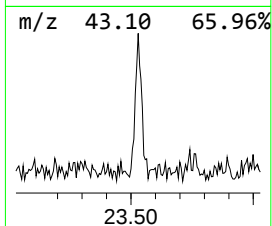
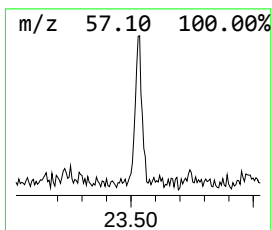
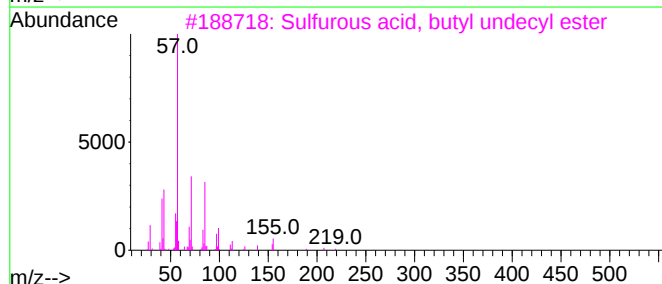
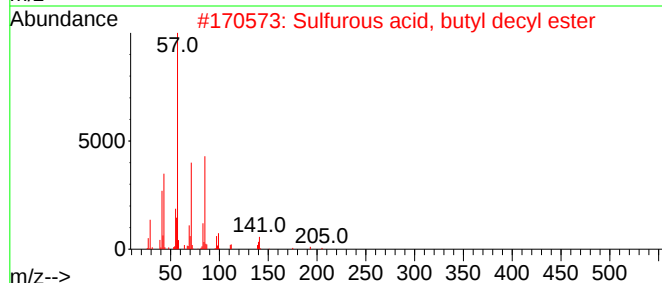
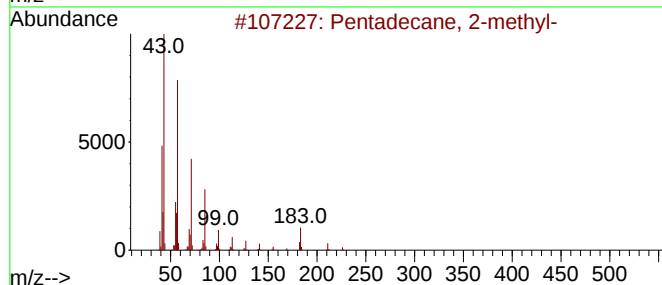
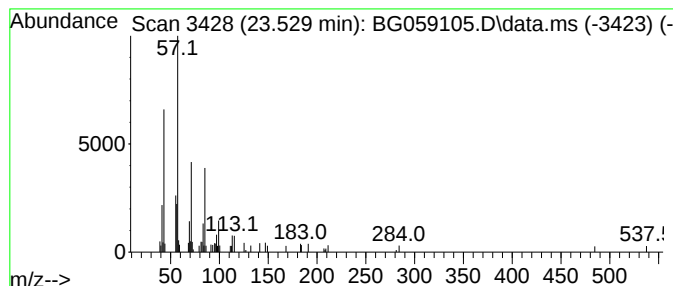
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.529	2.49 ng/ul	62102	Chrysene-d12	21.996

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane, 2-methyl-	226	C16H34	001560-93-6	78
2			Sulfurous acid, butyl decyl ester	278	C14H30O3S	1000309-17-7	72
3			Sulfurous acid, butyl undecyl ester	292	C15H32O3S	1000309-17-8	72
4			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	72
5			Undecane, 2,9-dimethyl-	184	C13H28	017301-26-7	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059105.D
Acq On : 20 Sep 2023 12:01
Operator : MA/JU
Sample : 04378-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BBJG5

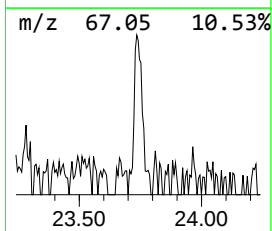
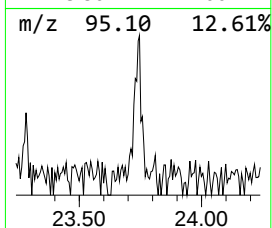
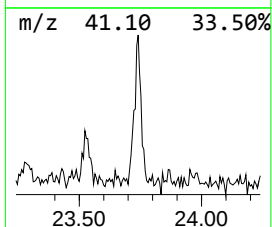
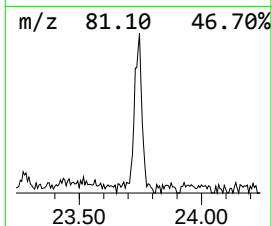
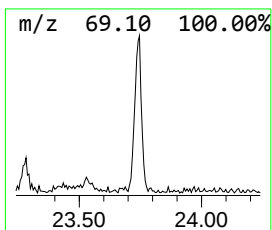
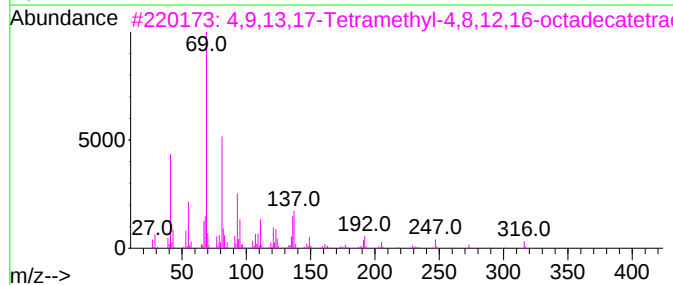
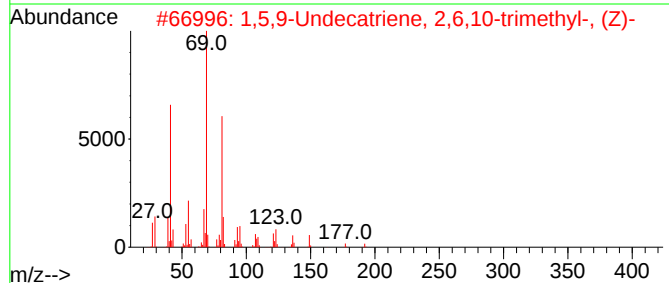
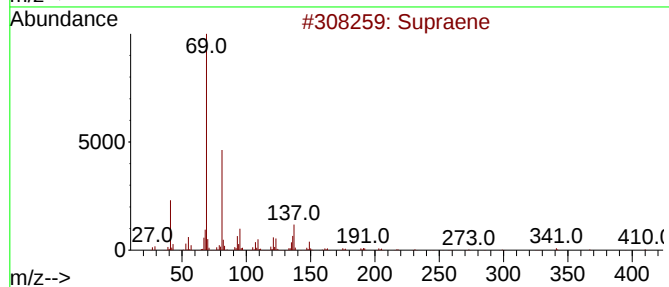
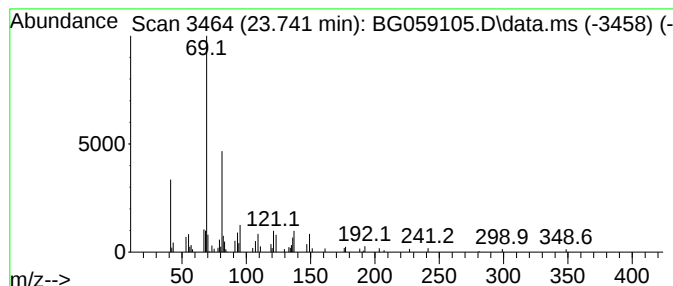
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Supraene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.741	3.72 ng/ul	92718	Chrysene-d12	21.996

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	80
2			1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	80
3			4,9,13,17-Tetramethyl-4,8,12,16-...	316	C22H36O	056882-09-8	78
4			6,10,14-Hexadecatrien-1-ol, 3,7,...	292	C20H36O	036237-66-8	64
5			1,6,10,14-Hexadecatetraen-3-ol, ...	290	C20H34O	001113-21-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059105.D
Acq On : 20 Sep 2023 12:01
Operator : MA/JU
Sample : 04378-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

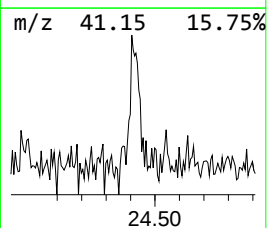
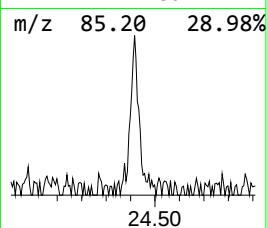
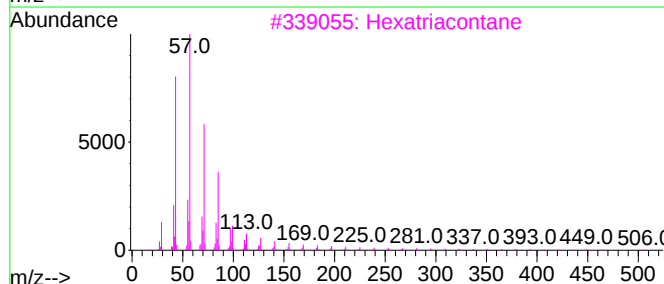
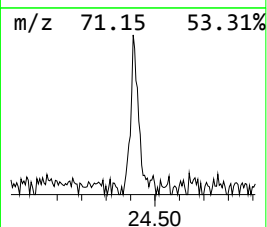
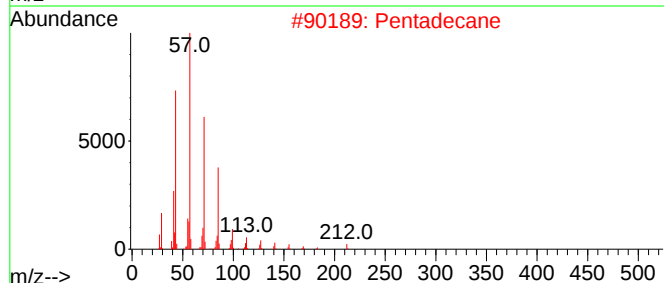
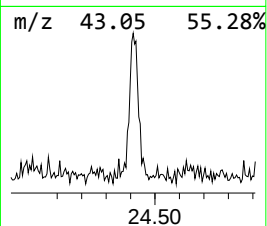
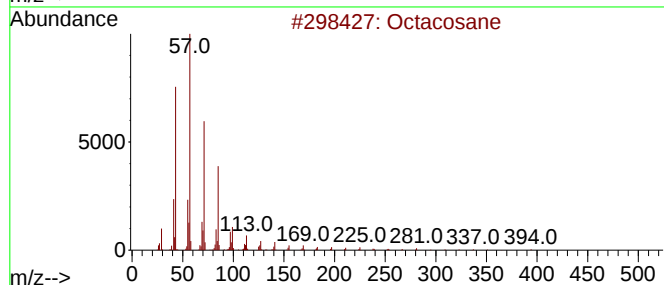
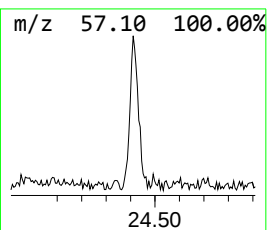
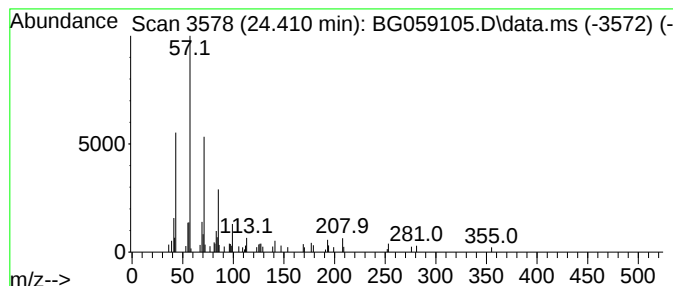
Instrument :
BNA_G
ClientSampleId :
BBJG5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.410	2.37 ng/ul	66310	Perylene-d12	25.544	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	74
2	Pentadecane	212	C15H32	000629-62-9	72
3	Hexatriacontane	507	C36H74	000630-06-8	72
4	Tetratetracontane	619	C44H90	007098-22-8	72
5	Heptacosane	380	C27H56	000593-49-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059105.D
Acq On : 20 Sep 2023 12:01
Operator : MA/JU
Sample : 04378-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BBJG5

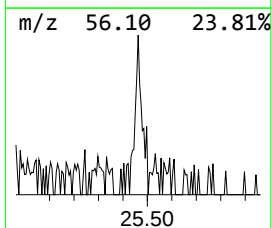
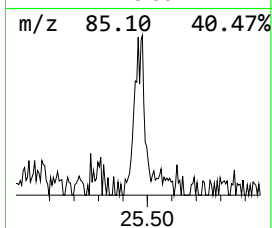
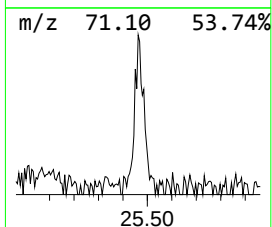
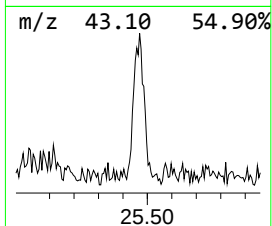
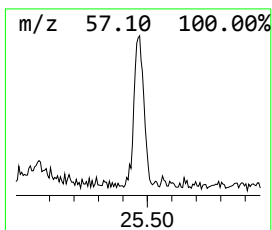
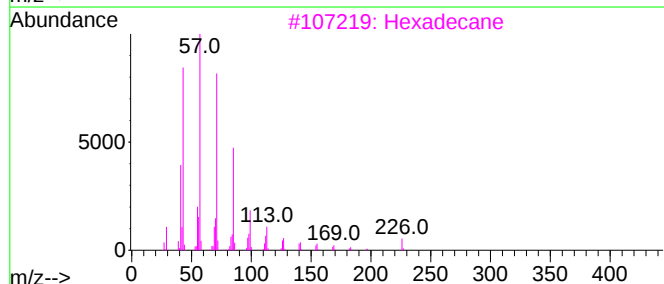
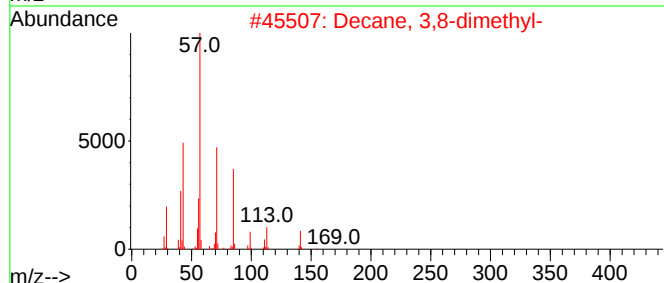
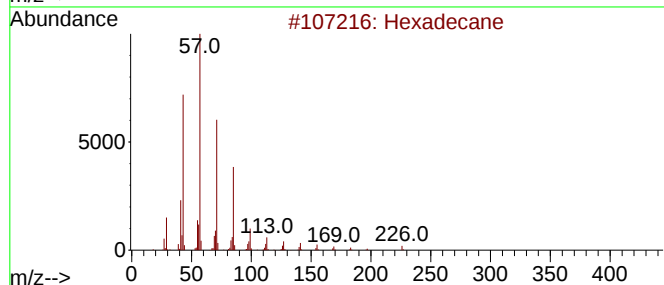
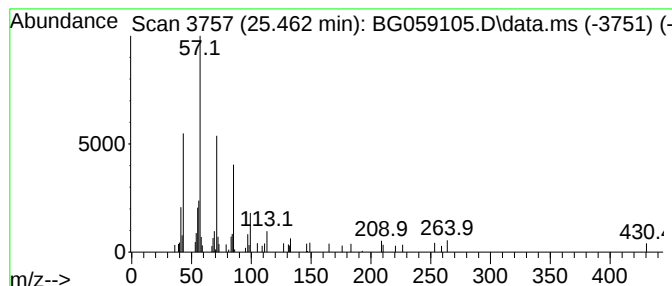
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.462	2.17 ng/ul	60652	Perylene-d12	25.544

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	72
2		Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	72
3		Hexadecane	226	C16H34	000544-76-3	64
4		Tetracosane	338	C24H50	000646-31-1	64
5		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059105.D
Acq On : 20 Sep 2023 12:01
Operator : MA/JU
Sample : 04378-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BBJG5

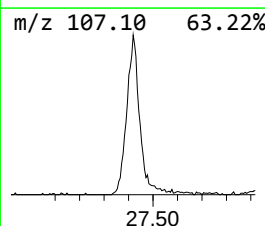
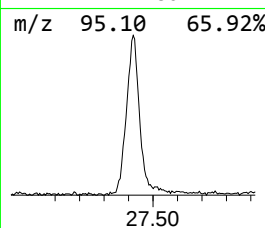
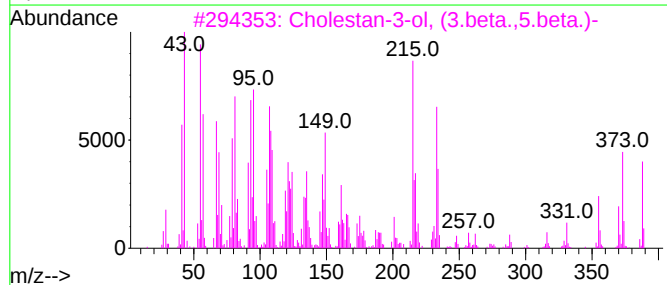
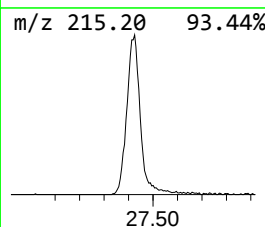
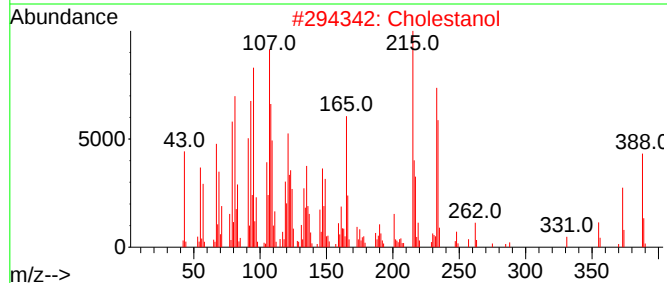
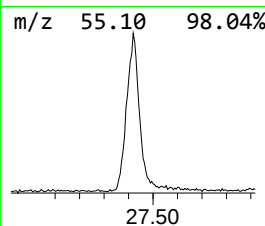
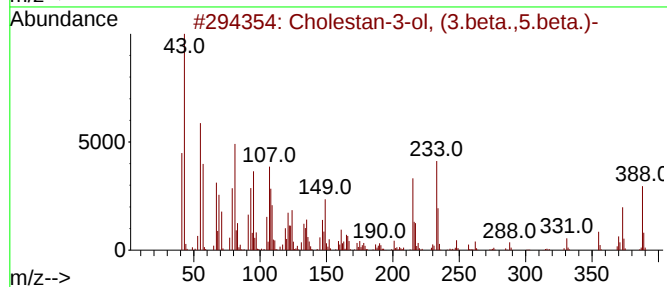
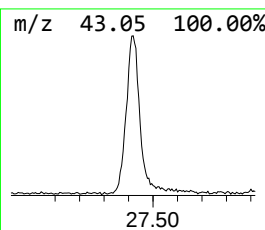
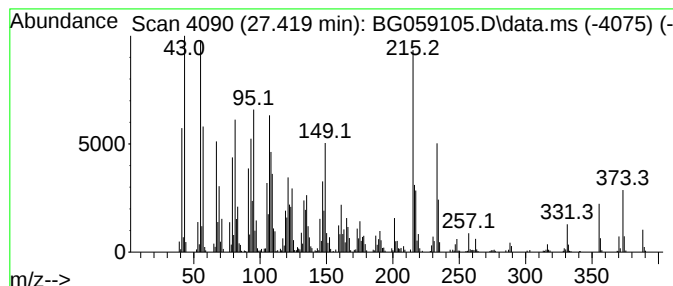
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Cholestan-3-ol, (3.beta.,5.... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.419	101.26 ng/ul	2835050	Perylene-d12	25.544

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cholestan-3-ol, (3.beta.,5.beta.)-	388	C27H48O	000360-68-9	68
2			Cholestanol	388	C27H48O	000080-97-7	64
3			Cholestan-3-ol, (3.beta.,5.beta.)-	388	C27H48O	000360-68-9	59
4			Cholestan-3-ol	388	C27H48O	027409-41-2	50
5			cis-3-[2-Hydroxy-4-(2-methyl-2-n...	332	C22H36O2	1000370-41-9	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059105.D
Acq On : 20 Sep 2023 12:01
Operator : MA/JU
Sample : 04378-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BBJG5

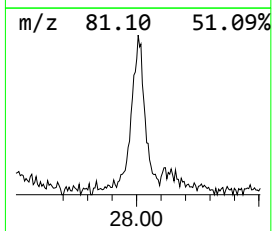
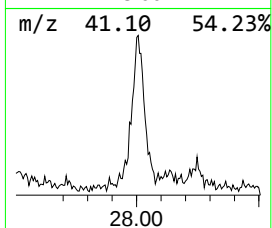
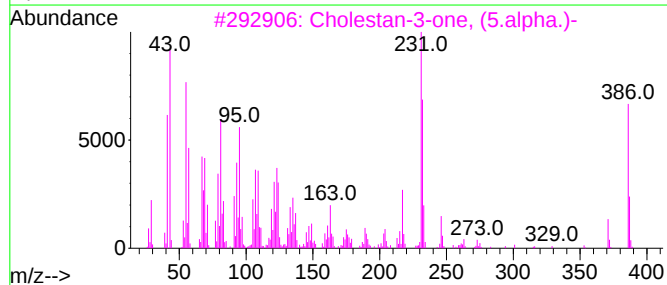
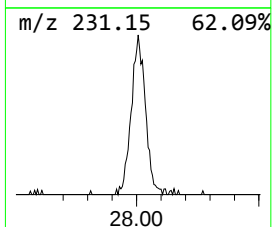
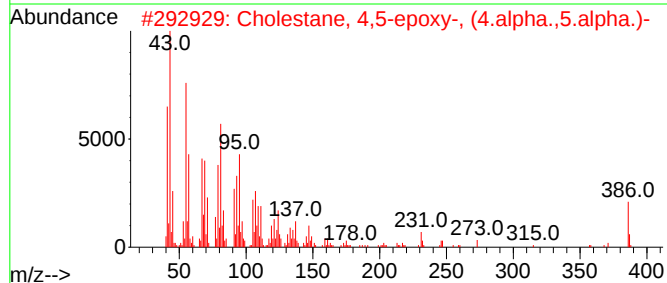
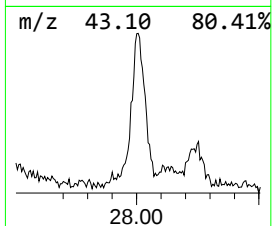
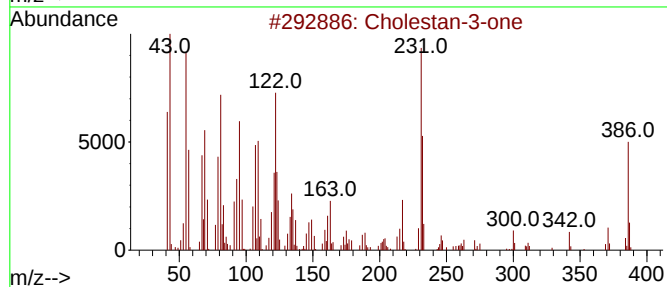
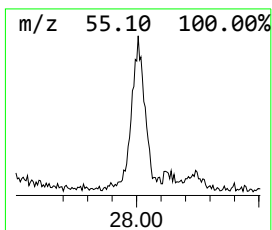
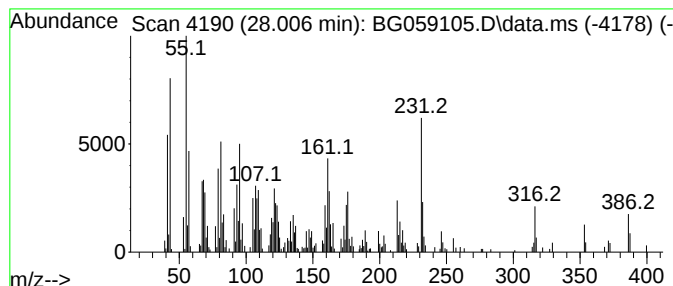
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 Cholestan-3-one Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.006	26.44 ng/ul	740267	Perylene-d12	25.544

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cholestan-3-one	386	C27H46O	015600-08-5	62
2			Cholestane, 4,5-epoxy-, (4.alpha...	386	C27H46O	006079-19-2	56
3			Cholestan-3-one, (5.alpha.)-	386	C27H46O	000566-88-1	47
4			Cholestan-2-one	386	C27H46O	1010210-43-4	38
5			Cholestan-3-one, (5.beta.)-	386	C27H46O	000601-53-6	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059105.D
Acq On : 20 Sep 2023 12:01
Operator : MA/JU
Sample : 04378-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BBJG5

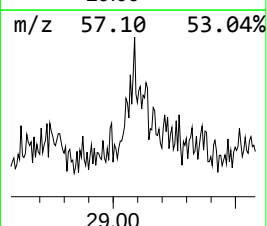
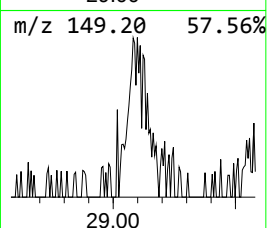
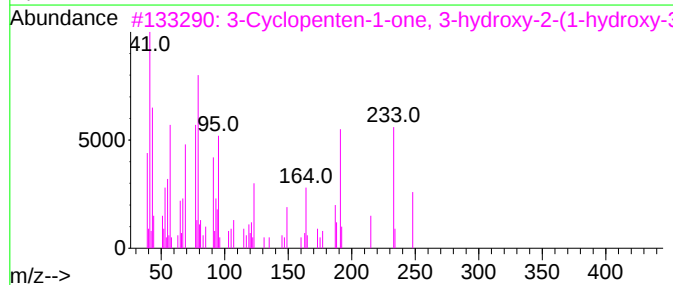
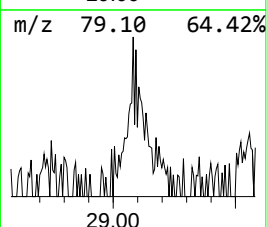
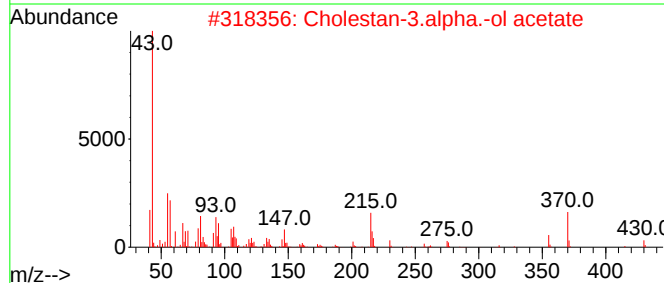
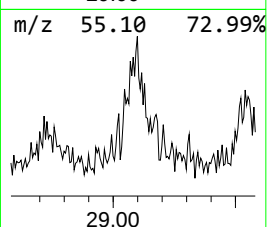
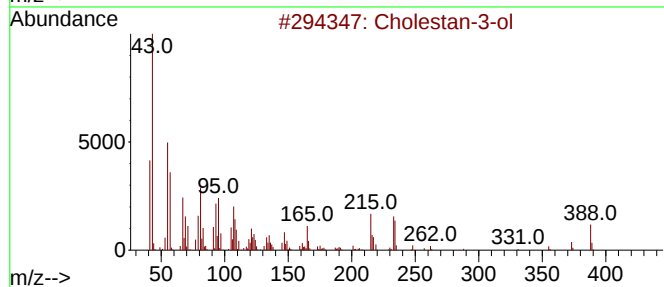
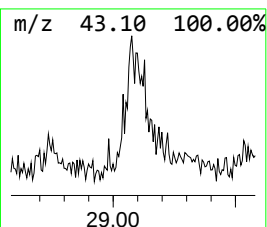
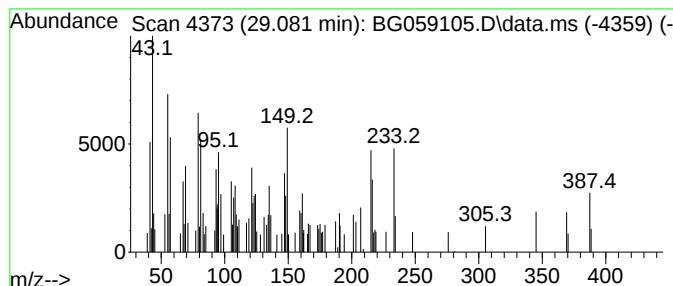
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.081	3.73 ng/ul	104382	Perylene-d12	25.544

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cholestan-3-ol	388	C27H48O	027409-41-2	43
2			Cholestan-3.alpha.-ol acetate	430	C29H50O2	022213-55-4	14
3			3-Cyclopenten-1-one, 3-hydroxy-2...	248	C15H20O3	038574-23-1	12
4			5.beta.-Cholestan-3.alpha.-ol, 2...	458	C31H54O2	1000514-95-7	10
5			5.beta.-Cholestan-3.alpha.-ol, a...	430	C29H50O2	033157-47-0	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059105.D
Acq On : 20 Sep 2023 12:01
Operator : MA/JU
Sample : 04378-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BBJG5

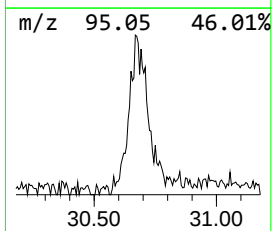
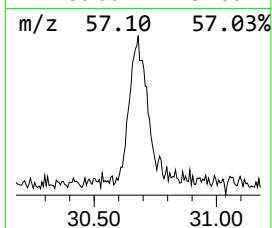
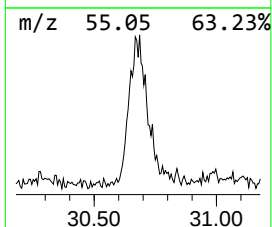
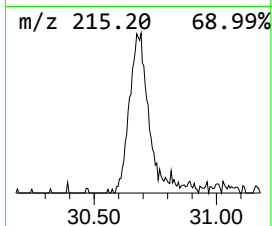
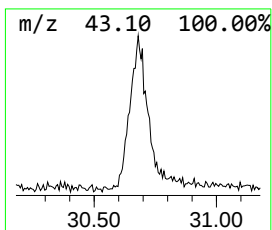
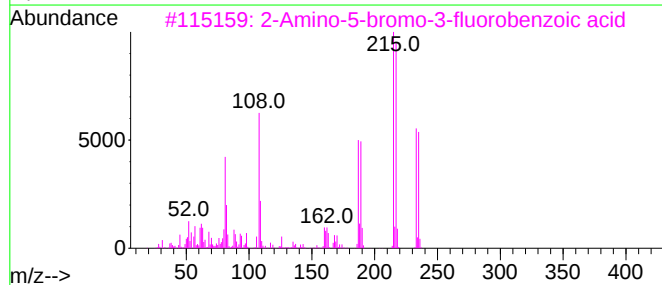
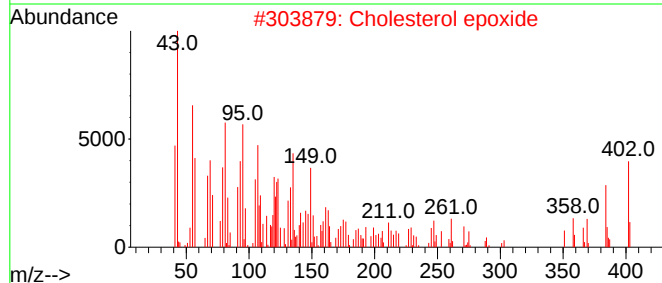
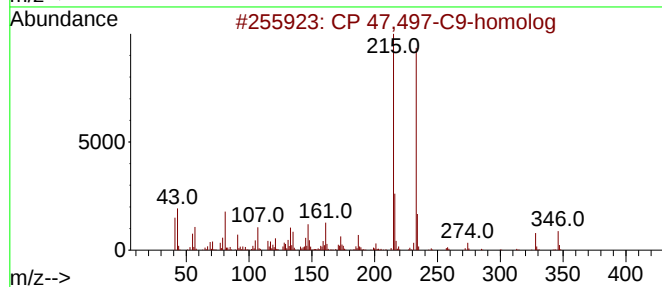
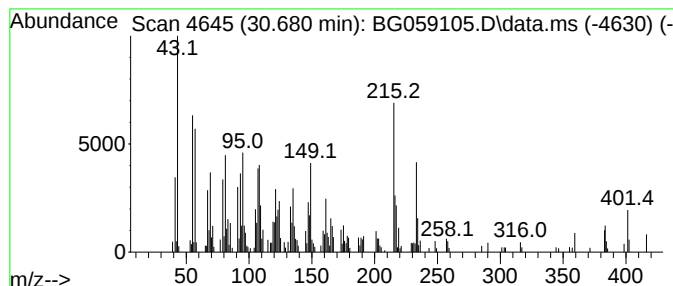
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-02 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.680	27.90 ng/ul	781154	Perylene-d12	25.544

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	CP	47,497	C9-homolog	346	C23H38O2	070435-08-4	49
2	Cholesterol	epoxide		402	C27H46O2	001250-95-9	35
3	2-Amino-5-bromo-3-fluorobenzoic	...		233	C7H5BrFNO2	874784-14-2	27
4	5-(2-Methyloctan-2-yl)-3-(1R,3R-...			332	C22H36O2	070434-92-3	27
5	CP	47,497	(n=7)	318	C21H34O2	070434-82-1	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059105.D
Acq On : 20 Sep 2023 12:01
Operator : MA/JU
Sample : 04378-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BBJG5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Tetrachloroethy...	4.916	7.8	ng/ul	72150	1	8.300	184937	20.0
n-Hexadecanoic ...	18.494	2.5	ng/ul	66185	4	17.677	532863	20.0
(DEL) Alkane: S...	23.529	2.5	ng/ul	62102	5	21.996	498249	20.0
Supraene	23.741	3.7	ng/ul	92718	5	21.996	498249	20.0
(DEL) Alkane: S...	24.410	2.4	ng/ul	66310	6	25.544	559940	20.0
(DEL) Alkane: S...	25.462	2.2	ng/ul	60652	6	25.544	559940	20.0
Cholestan-3-ol,...	27.419	101.3	ng/ul	2835050	6	25.544	559940	20.0
Cholestan-3-one	28.006	26.4	ng/ul	740267	6	25.544	559940	20.0
unknown-01	29.081	3.7	ng/ul	104382	6	25.544	559940	20.0
unknown-02	30.680	27.9	ng/ul	781154	6	25.544	559940	20.0