LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100923\
Data File : BG059408.D
Acq On : 9 Oct 2023 22:42
Operator : MA/JU
Sample : 04744-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BH580

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-BG100723.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BG059408.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.967	93	98	108	rVB	57611	109298	4.38%	0.674%
2	4.037	108	110	111	rBV	8852	7476	0.30%	0.046%
3	4.836	239	246	253	rBV	1632866	2493978	100.00%	15.381%
4	5.136	292	297	302	rVB2	142952	208123	8.35%	1.284%
5	5.306	322	326	329	rVV3	13574	16579	0.66%	0.102%
6	5.359	330	335	342	rVB2	73391	113780	4.56%	0.702%
7	5.688	387	391	393	rBV	5659	7517	0.30%	0.046%
8	5.788	405	408	411	rBV	6477	6844	0.27%	0.042%
9	5.870	419	422	426	rVB	4972	6929	0.28%	0.043%
10	6.411	512	514	516	rBV	7942	6542	0.26%	0.040%
11	6.446	516	520	523	rVV3	15718	20945	0.84%	0.129%
12	7.122	631	635	638	rBV	7637	14208	0.57%	0.088%
13	7.351	670	674	675	rBV3	9559	13148	0.53%	0.081%
14	7.545	702	707	714	rVB	192350	319416	12.81%	1.970%
15	7.633	718	722	727	rVB2	26541	41740	1.67%	0.257%
16	7.745	729	741	749	rBV	247984	481692	19.31%	2.971%
17	8.226	817	823	831	rBV3	134554	238721	9.57%	1.472%
18	8.297	831	835	843	rVB2	23881	39388	1.58%	0.243%
19	8.714	905	906	910	rBV	8424	7674	0.31%	0.047%
20	8.920	935	941	949	rBV3	34396	65752	2.64%	0.405%
21	9.072	964	967	970	rBV	6094	9326	0.37%	0.058%
22	9.102	970	972	976	rVV2	13499	18230	0.73%	0.112%
23	9.425	1015	1027	1034	rBV	234167	468391	18.78%	2.889%
24	9.501	1034	1040	1047	rVB	160465	279579	11.21%	1.724%
25	9.789	1086	1089	1091	rBV	5258	7522	0.30%	0.046%
26	9.872	1101	1103	1107	rVB	7100	6493	0.26%	0.040%
27	9.936	1110	1114	1116	rBV	6252	6969	0.28%	0.043%
28	10.142	1142	1149	1157	rBV2	187125	347828	13.95%	2.145%
29	10.265	1164	1170	1171	rBV2	12000	18839	0.76%	0.116%
30	10.671	1232	1239	1248	rVB	267832	483183	19.37%	2.980%
31	10.847	1262	1269	1274	rBV5	51441	101758	4.08%	0.628%
32	10.906	1275	1279	1289	rVB3	65508	128479	5.15%	0.792%
33	11.070	1300	1307	1314	rVB	209337	383001	15.36%	2.362%
34	11.605	1396	1398	1400	rBV	5906	6613	0.27%	0.041%
35	11.787	1425	1429	1435	rBV3	21305	39038	1.57%	0.241%
36	12.727	1584	1589	1594	rBV2	9846	17184	0.69%	0.106%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100923\
Data File : BG059408.D
Acq On : 9 Oct 2023 22:42
Operator : MA/JU
Sample : 04744-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BH580

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-BG100723.MA.M
Title : SVOA CALIBRATION

37	13.273	1679	1682	1687	rBV3	11829	19554	0.78%	0.121%
38	13.585	1732	1735	1736	rBV	5504	6999	0.28%	0.043%
39	13.679	1748	1751	1753	rBV3	9506	12854	0.52%	0.079%
40	14.014	1805	1808	1809	rBV	7092	8111	0.33%	0.050%
41	14.255	1843	1849	1855	rVB	586143	859359	34.46%	5.300%
42	14.496	1888	1890	1894	rBV	5954	6873	0.28%	0.042%
43	14.560	1895	1901	1908	rVV	332096	519171	20.82%	3.202%
44	14.860	1946	1952	1960	rVB	368732	574286	23.03%	3.542%
45	14.983	1971	1973	1976	rBV2	8799	8174	0.33%	0.050%
46	15.054	1980	1985	1990	rBV3	43737	71415	2.86%	0.440%
47	15.724	2096	2099	2103	rBV	6403	11095	0.44%	0.068%
48	15.794	2108	2111	2113	rBV2	8903	8685	0.35%	0.054%
49	15.847	2114	2120	2126	rVV	898631	1338812	53.68%	8.257%
50	15.959	2133	2139	2145	rBV	338338	481272	19.30%	2.968%
51	16.552	2237	2240	2243	rBV2	12771	16373	0.66%	0.101%
52	16.793	2278	2281	2283	rBV	8384	9355	0.38%	0.058%
53	17.251	2354	2359	2365	rBV4	43621	69138	2.77%	0.426%
54	17.474	2395	2397	2403	rVB4	16199	24458	0.98%	0.151%
55	17.610	2414	2420	2427	rVB	459702	697114	27.95%	4.299%
56	17.709	2430	2437	2446	rVV2	640740	1005454	40.32%	6.201%
57	17.845	2455	2460	2471	rVB	139067	238914	9.58%	1.473%
58	18.432	2556	2560	2564	rVB4	35254	54833	2.20%	0.338%
59	18.984	2650	2654	2661	rVB2	48604	67916	2.72%	0.419%
60	19.825	2795	2797	2800	rVB3	16294	17209	0.69%	0.106%
61	19.983	2818	2824	2832	rVB	1083397	1457739	58.45%	8.990%
62	21.211	3029	3033	3039	rVB	117573	177598	7.12%	1.095%
63	21.910	3146	3152	3159	rBV	417566	693033	27.79%	4.274%
64	25.148	3696	3703	3714	rVB2	428718	1217145	48.80%	7.506%

Sum of corrected areas: 16215122

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100923\
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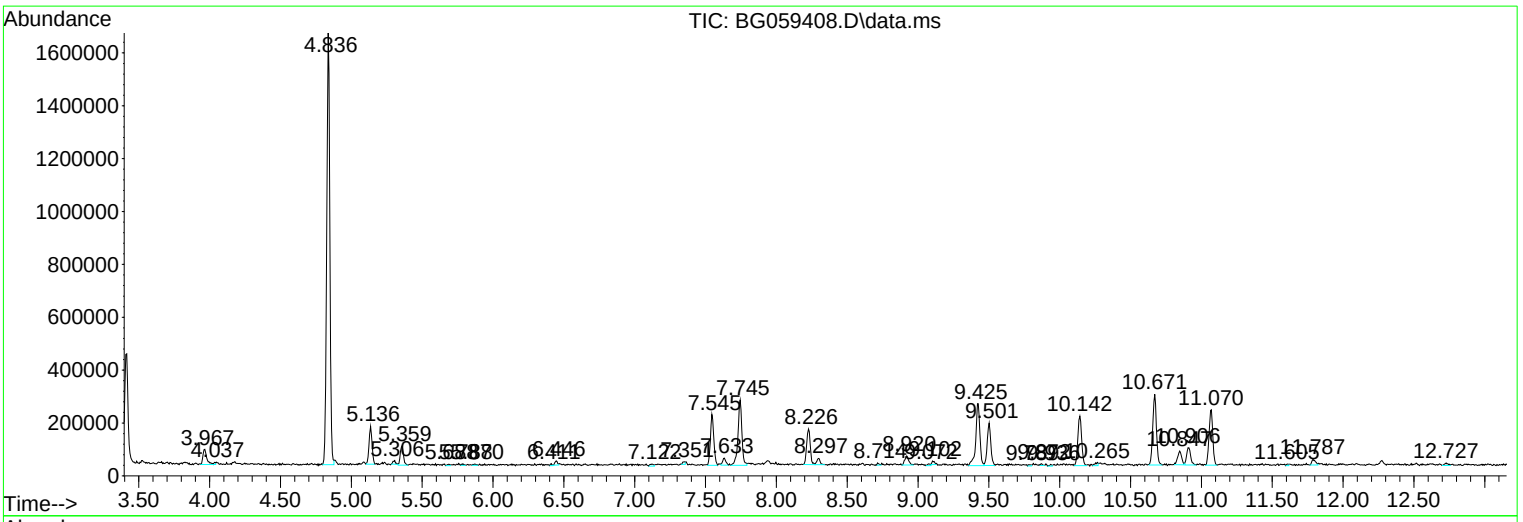
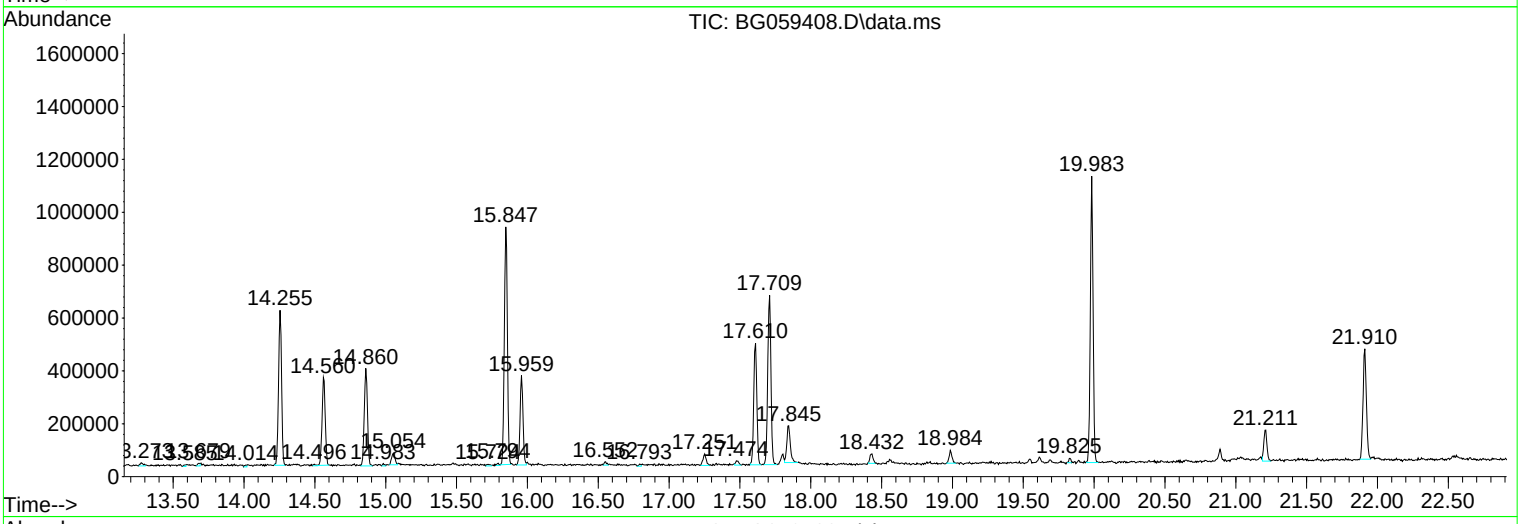
Misc :
ALS Vial : 18 Sample Multiplier: 1

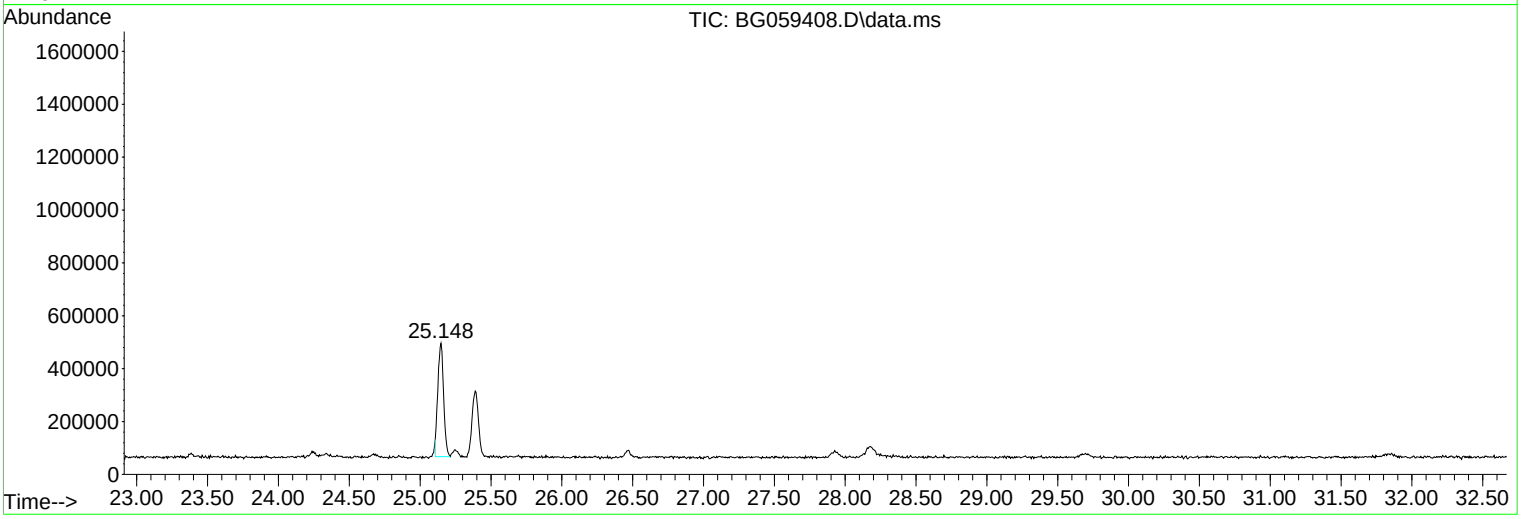
Instrument :
BNA_G
ClientSampleId :
BH580

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
Quant Title : SVOA CALIBRATION

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







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Data File : BG059408.D
Acq On : 9 Oct 2023 22:42
Operator : MA/JU
Sample : 04744-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH580

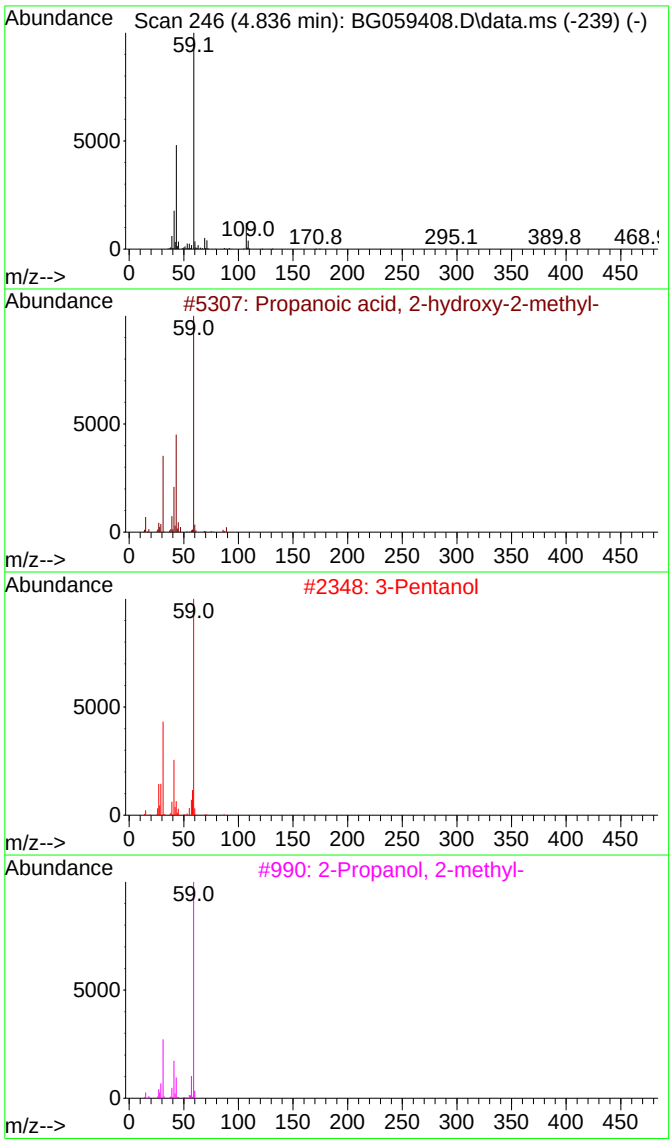
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
Quant Title : SVOA CALIBRATION

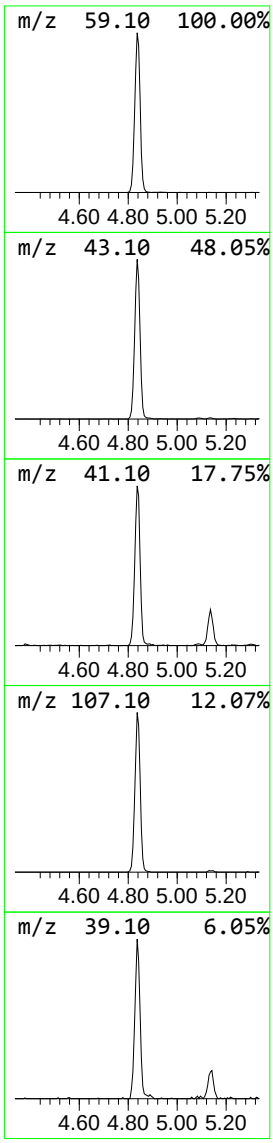
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.836	208.94 ng/ul	2493980	1,4-Dichlorobenzene-d4	8.226

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	45
2	3-Pentanol	88	C5H12O	000584-02-1	42
3	2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	40
4	3-Pentanol	88	C5H12O	000584-02-1	36
5	3-Pentanol	88	C5H12O	000584-02-1	36





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100923\
Data File : BG059408.D
Acq On : 9 Oct 2023 22:42
Operator : MA/JU
Sample : 04744-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH580

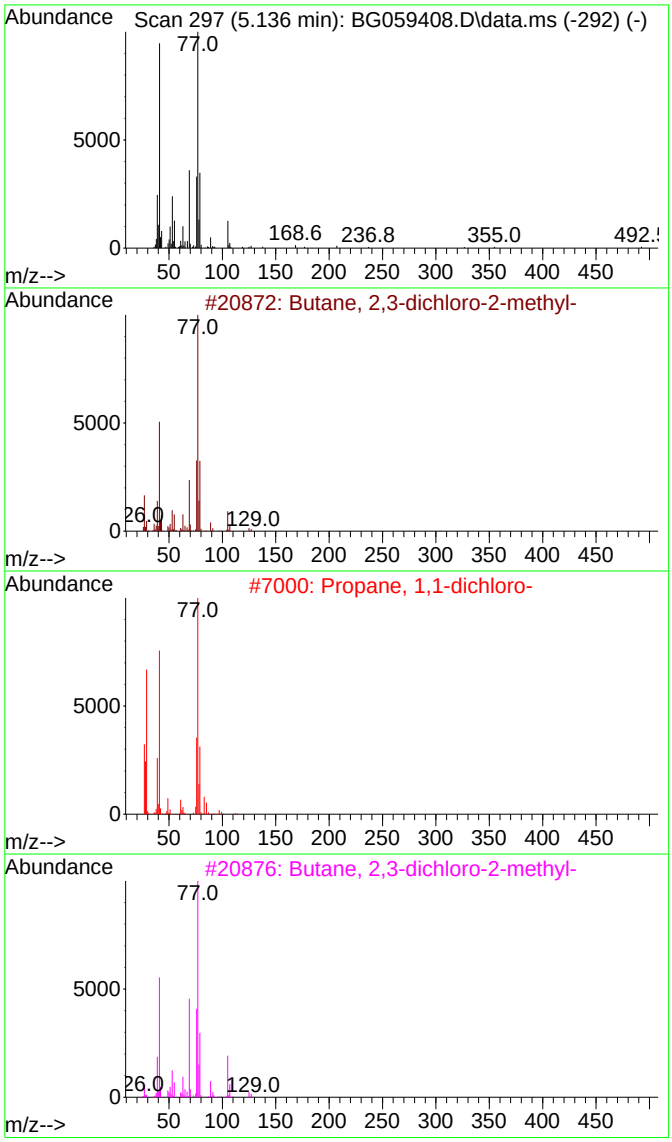
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
Quant Title : SVOA CALIBRATION

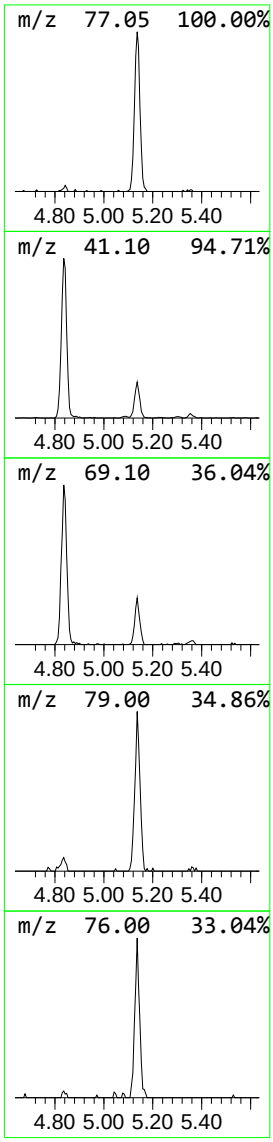
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2,3-dichloro-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.136	17.44 ng/ul	208123	1,4-Dichlorobenzene-d4	8.226

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dichloro-2-methyl-	140	C5H10Cl2	000507-45-9	72
2	Propane, 1,1-dichloro-	112	C3H6Cl2	000078-99-9	42
3	Butane, 2,3-dichloro-2-methyl-	140	C5H10Cl2	000507-45-9	38
4	Propane, 2-bromo-1-chloro-	156	C3H6BrCl	003017-95-6	25
5	Propane, 2-bromo-1-chloro-	156	C3H6BrCl	003017-95-6	23





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100923\
Data File : BG059408.D
Acq On : 9 Oct 2023 22:42
Operator : MA/JU
Sample : 04744-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH580

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
Quant Title : SVOA CALIBRATION

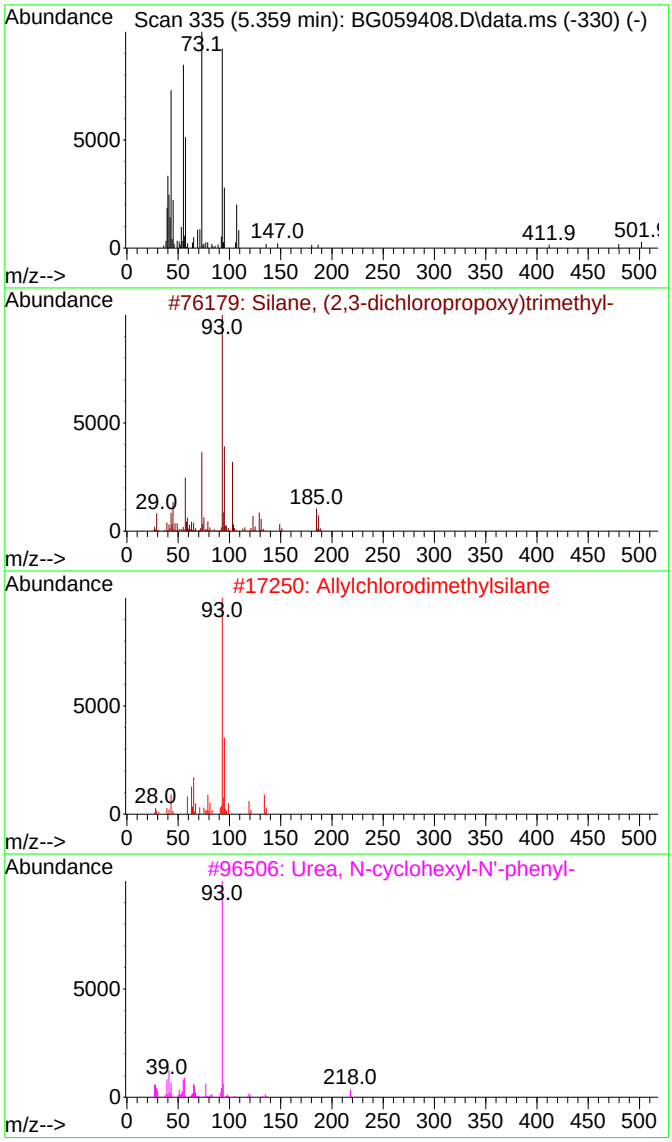
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

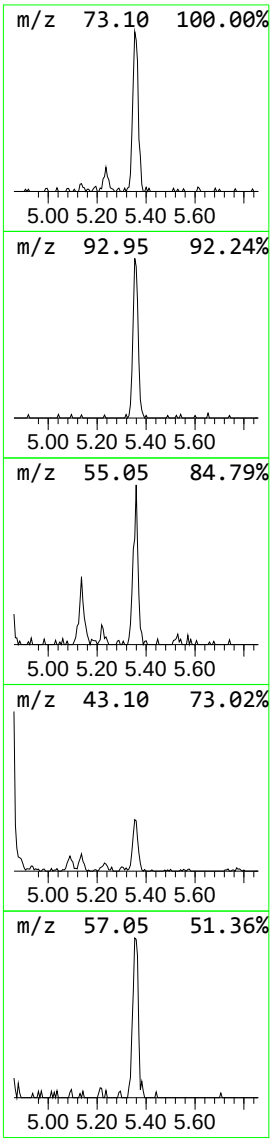
****

Peak Number 3 unknown-02 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.359	9.53 ng/ul	113780	1,4-Dichlorobenzene-d4	8.226

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Silane, (2,3-dichloropropoxy)tri...	200	C6H14Cl2OSi	018151-22-9	33
2	Allylchlorodimethylsilane	134	C5H11ClSi	004028-23-3	32
3	Urea, N-cyclohexyl-N'-phenyl-	218	C13H18N2O	000886-59-9	25
4	Propanamide, 3-bromo-N-phenyl-	227	C9H10BrNO	007661-07-6	17
5	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	17





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100923\
Data File : BG059408.D
Acq On : 9 Oct 2023 22:42
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ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH580

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
Quant Title : SVOA CALIBRATION

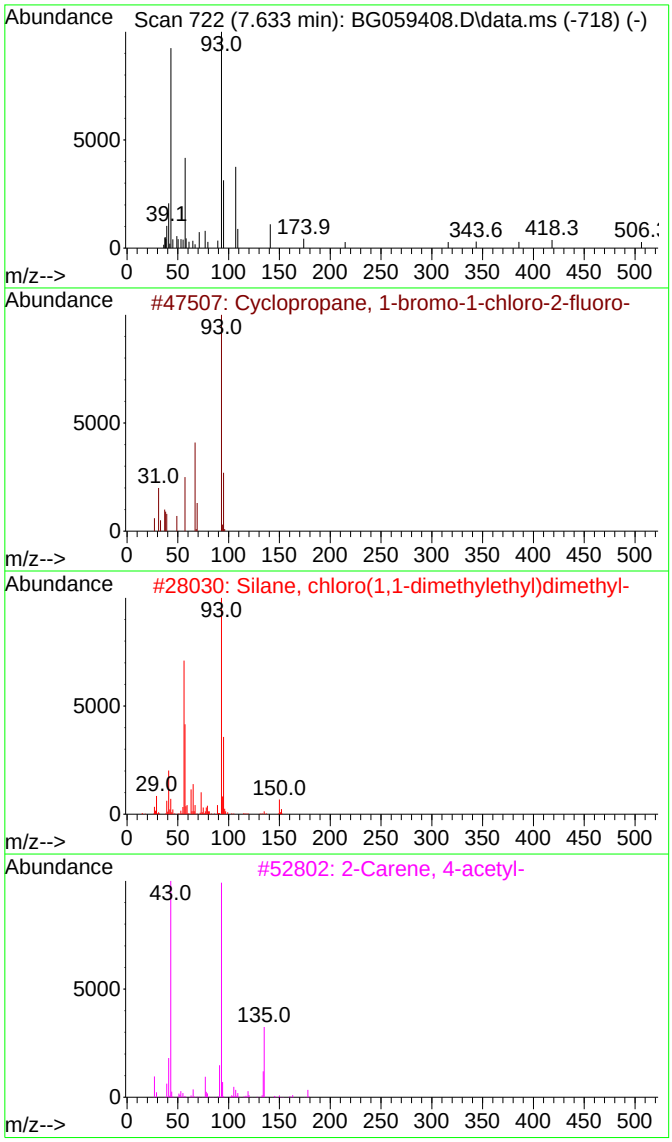
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

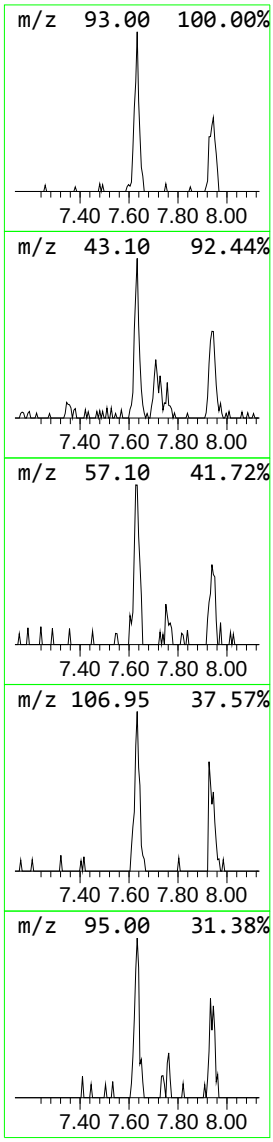
****

Peak Number 4 unknown-03 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.633	3.50 ng/ul	41740	1,4-Dichlorobenzene-d4	8.226

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1-bromo-1-chloro-2...	172	C3H3BrClF	024071-59-8	38
2	Silane, chloro(1,1-dimethylethyl...	150	C6H15ClSi	018162-48-6	9
3	2-Carene, 4-acetyl-	178	C12H18O	1000163-28-1	9
4	2-Butanol, 1,4-dichloro-	142	C4H8Cl2O	002419-74-1	9
5	2-Furancarboxaldehyde, 5-nitro-	141	C5H3NO4	000698-63-5	7





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100923\
Data File : BG059408.D
Acq On : 9 Oct 2023 22:42
Operator : MA/JU
Sample : 04744-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH580

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
Quant Title : SVOA CALIBRATION

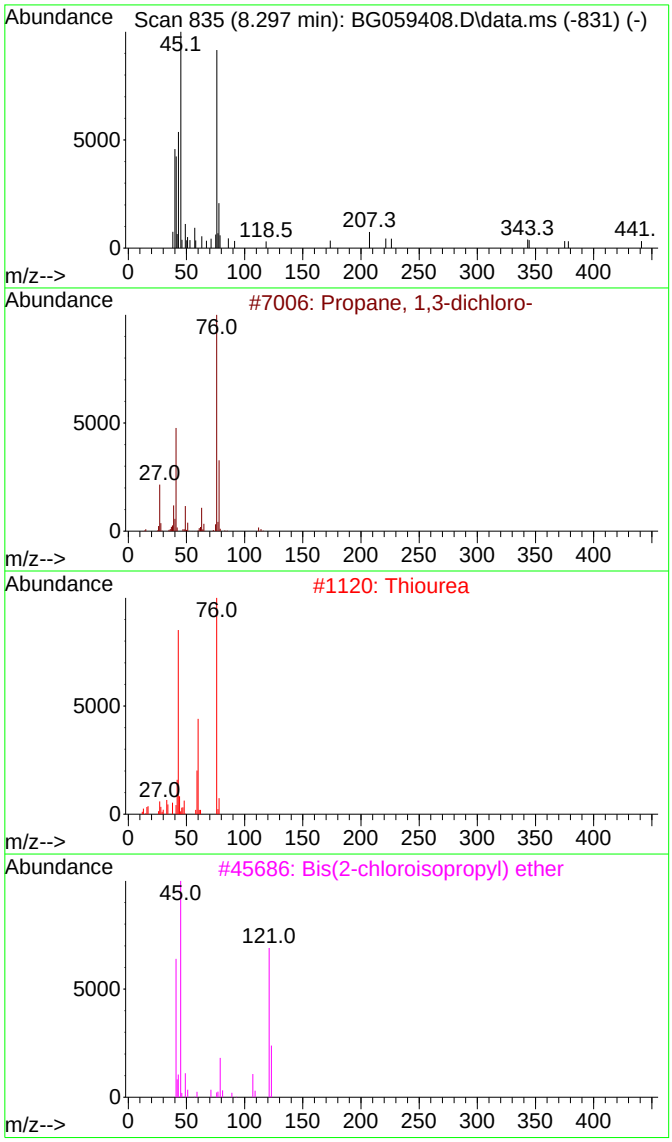
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

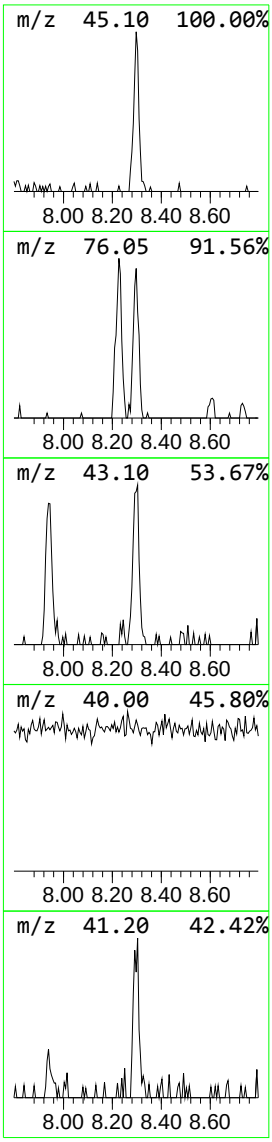
****

Peak Number 5 unknown-04 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.297	3.30 ng/ul	39388	1,4-Dichlorobenzene-d4	8.226

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 1,3-dichloro-	112	C3H6Cl2	000142-28-9	45
2	Thiourea	76	CH4N2S	000062-56-6	9
3	Bis(2-chloroisopropyl) ether	170	C6H12Cl2O	039638-32-9	9
4	Methane, chloromethoxy-	80	C2H5ClO	000107-30-2	9
5	2-Pentanol, 3-chloro-4-methyl-, ...	136	C6H13ClO	074685-47-5	9





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Data File : BG059408.D
Acq On : 9 Oct 2023 22:42
Operator : MA/JU
Sample : 04744-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH580

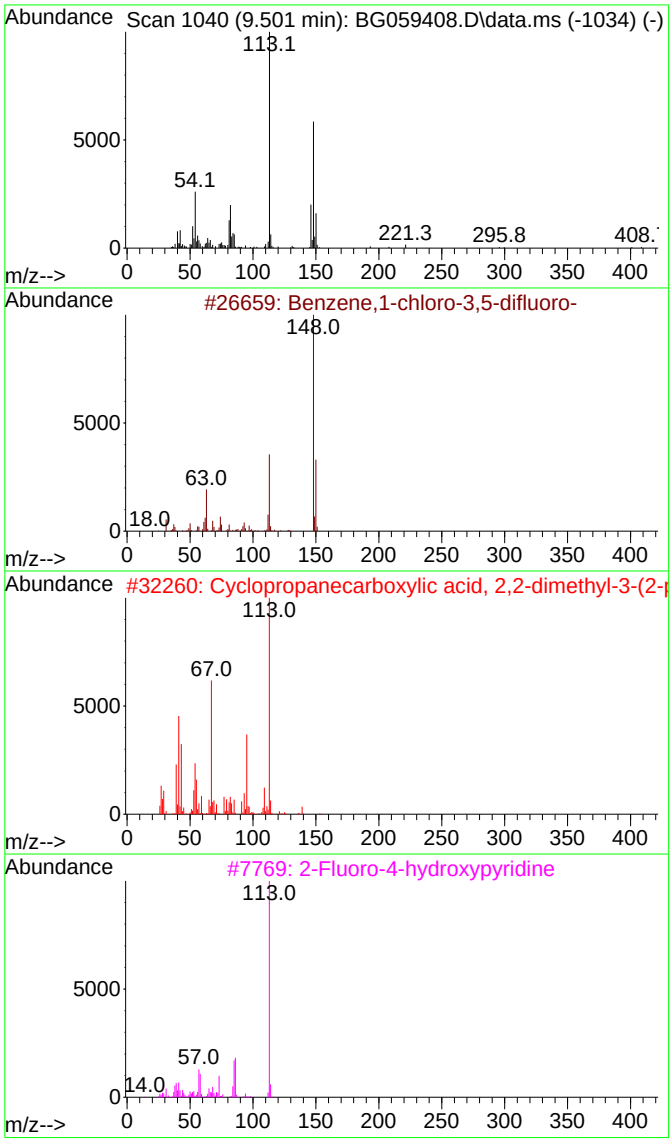
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
Quant Title : SVOA CALIBRATION

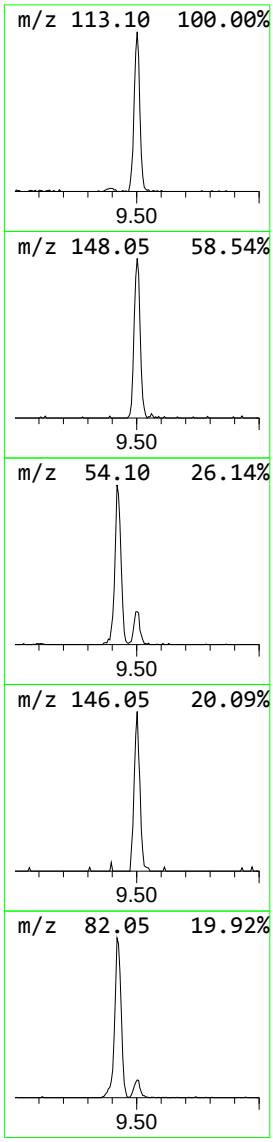
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene,1-chloro-3,5-difluoro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.501	23.42 ng/ul	279579	1,4-Dichlorobenzene-d4	8.226

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene,1-chloro-3,5-difluoro-	148	C6H3ClF2	001435-43-4	50
2	Cyclopropanecarboxylic acid, 2,2...	154	C9H14O2	074685-76-0	32
3	2-Fluoro-4-hydroxypyridine	113	C5H4FNO	022282-69-5	27
4	4-Methylamino-2(5H)-furanone	113	C5H7NO2	1000427-45-5	16
5	Ethylphoshonic acid, fluoroanhyd...	224	C10H22FO2P	1000273-53-4	16





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100923\
Data File : BG059408.D
Acq On : 9 Oct 2023 22:42
Operator : MA/JU
Sample : 04744-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH580

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
Quant Title : SVOA CALIBRATION

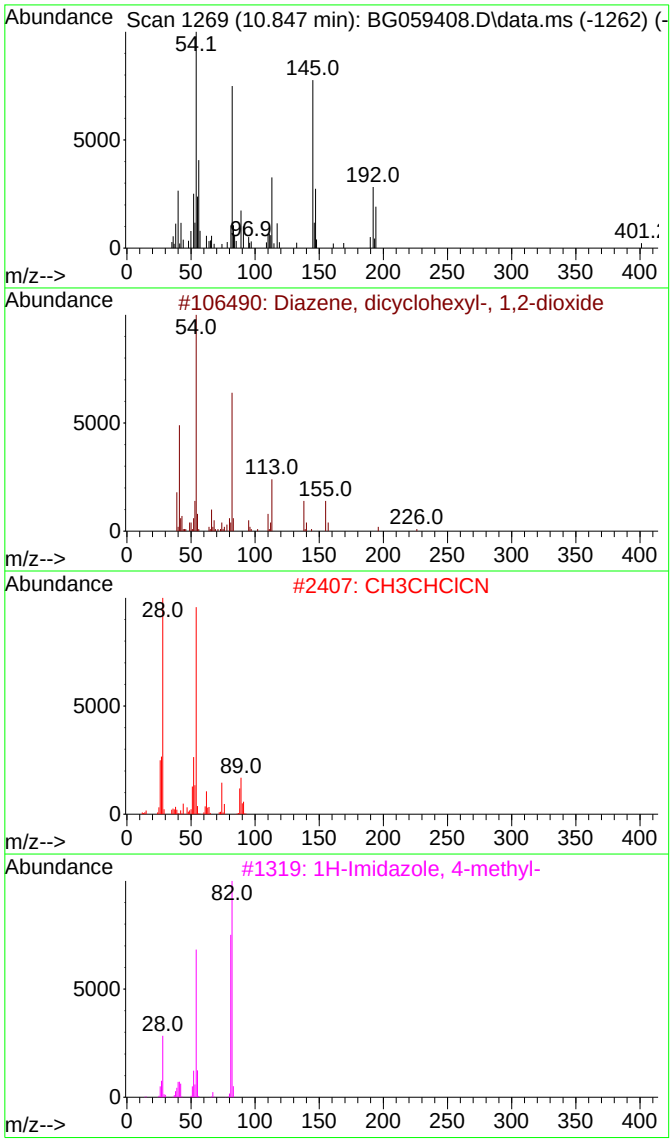
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

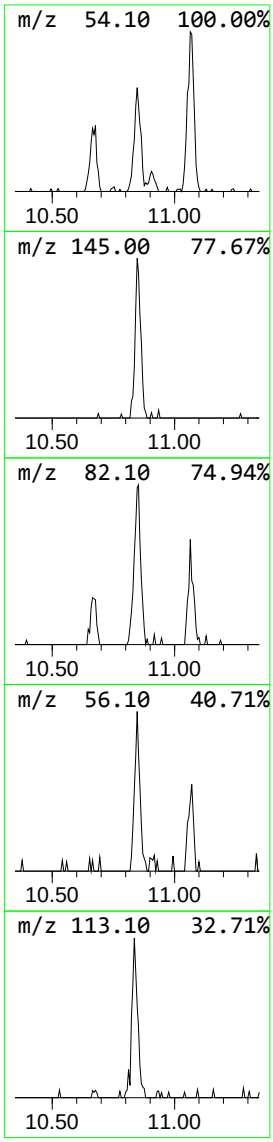
****

Peak Number 7 unknown-05 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.847	5.31 ng/ul	101758	Naphthalene-d8	11.064

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diazene, dicyclohexyl-, 1,2-dioxide	226	C12H22N2O2	003378-45-8	47
2	CH3CHClCN	89	C3H4ClN	001617-17-0	43
3	1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	42
4	6-Chloro-2,3-dihydroxypyridine	145	C5H4ClNO2	1000341-18-9	38
5	N-Methylmaleimide	111	C5H5NO2	000930-88-1	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100923\
Data File : BG059408.D
Acq On : 9 Oct 2023 22:42
Operator : MA/JU
Sample : 04744-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH580

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
Quant Title : SVOA CALIBRATION

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

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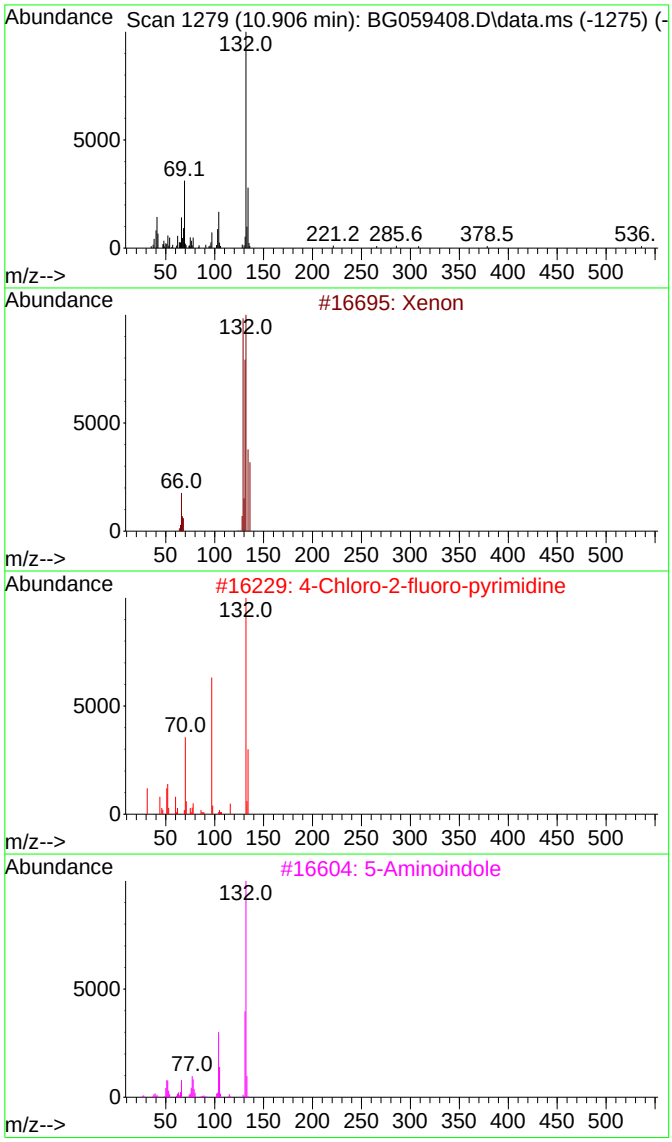
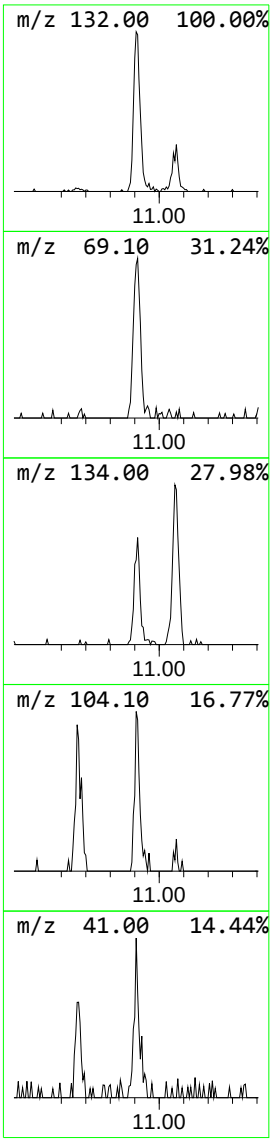
Peak Number 8 Xenon Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.906	6.71 ng/ul	128479	Naphthalene-d8	11.064

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Xenon	132	Xe	007440-63-3	59
2	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	59
3	5-Aminoindole	132	C8H8N2	005192-03-0	58
4	5-Aminoindole	132	C8H8N2	005192-03-0	58
5	1H-Indol-4-amine	132	C8H8N2	005192-23-4	53





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Data File : BG059408.D
Acq On : 9 Oct 2023 22:42
Operator : MA/JU
Sample : 04744-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH580

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
Quant Title : SVOA CALIBRATION

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

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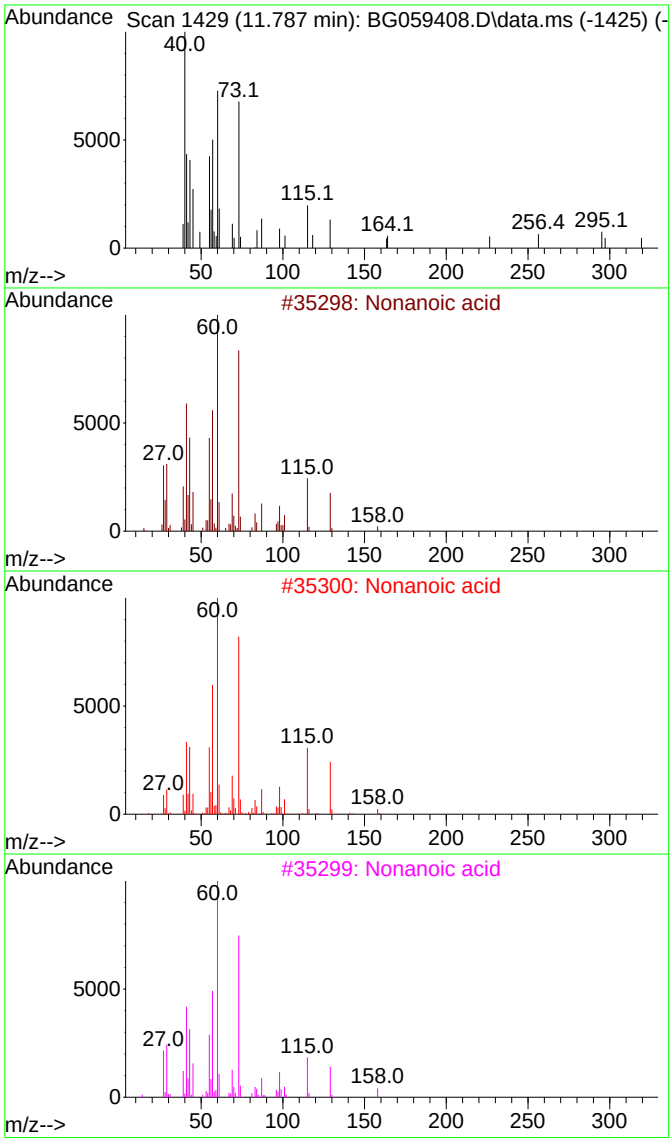
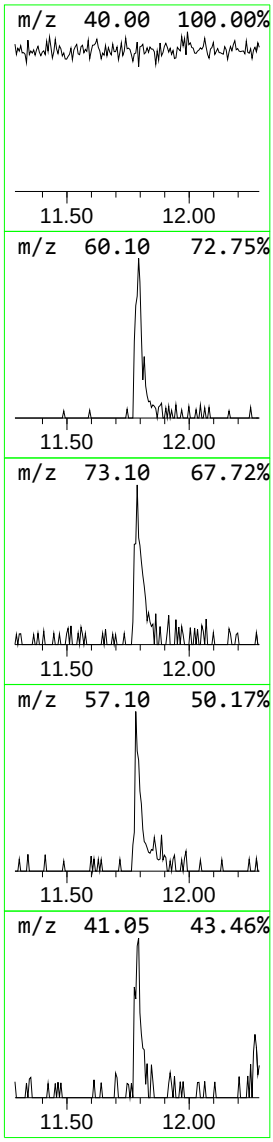
Peak Number 9 Nonanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.787	2.04 ng/ul	39038	Naphthalene-d8	11.064

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanoic acid	158	C9H18O2	000112-05-0	80
2	Nonanoic acid	158	C9H18O2	000112-05-0	72
3	Nonanoic acid	158	C9H18O2	000112-05-0	64
4	Nonanoic acid	158	C9H18O2	000112-05-0	64
5	Nonanoic acid	158	C9H18O2	000112-05-0	64





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Data File : BG059408.D
Acq On : 9 Oct 2023 22:42
Operator : MA/JU
Sample : 04744-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH580

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
Quant Title : SVOA CALIBRATION

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

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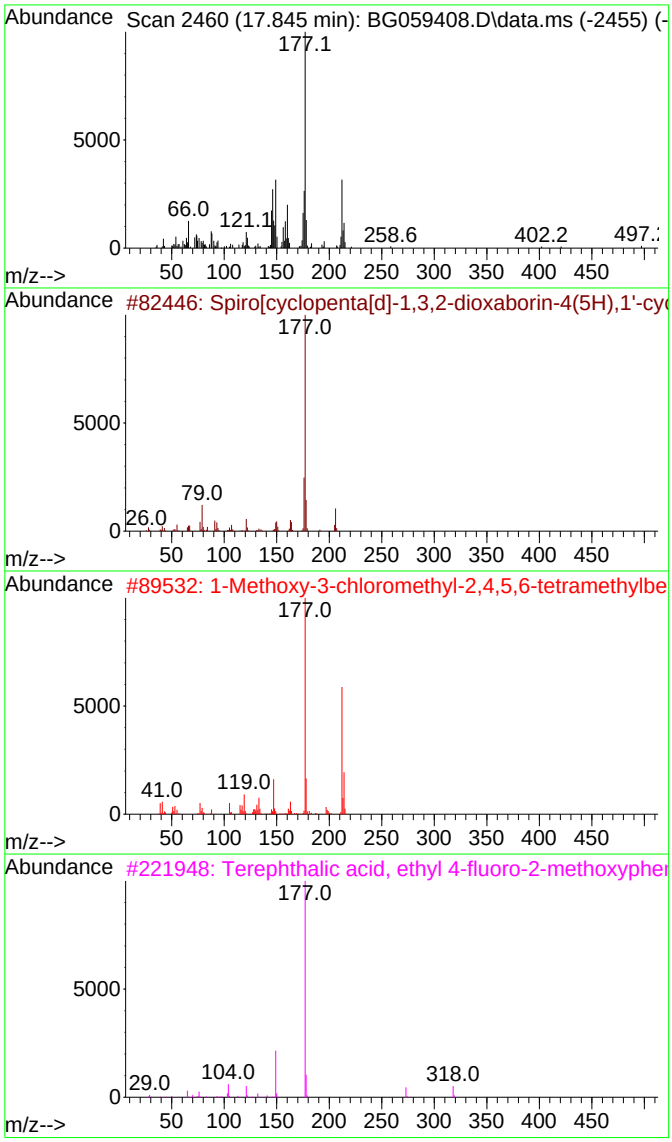
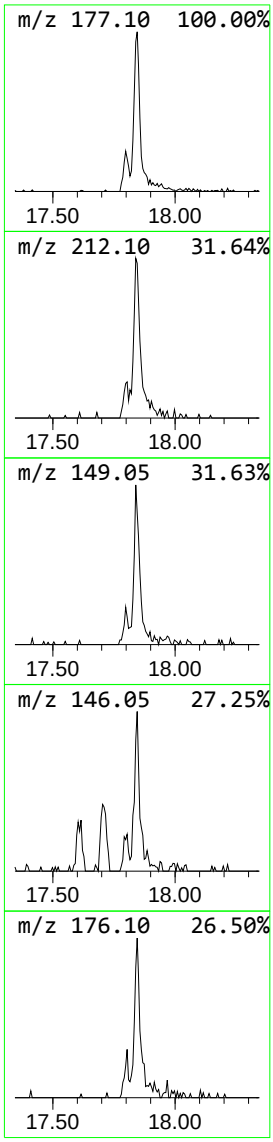
Peak Number 10 unknown-06 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.845	6.85 ng/ul	238914	Phenanthrene-d10	17.610

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Spiro[cyclopenta[d]-1,3,2-dioxab...	206	C12H19BO2	062238-26-0	43
2	1-Methoxy-3-chloromethyl-2,4,5,6...	212	C12H17ClO	073942-80-0	43
3	Terephthalic acid, ethyl 4-fluor...	318	C17H15FO5	1000415-82-7	38
4	Benzoic acid, 2,3,5,6-tetrafluor...	234	C10H6F4O2	1000460-51-7	38
5	Benzene, 1-chloro-2-(dichloromet...	212	C7H4Cl3F	062476-62-4	38





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Data File : BG059408.D
Acq On : 9 Oct 2023 22:42
Operator : MA/JU
Sample : 04744-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH580

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
Quant Title : SVOA CALIBRATION

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

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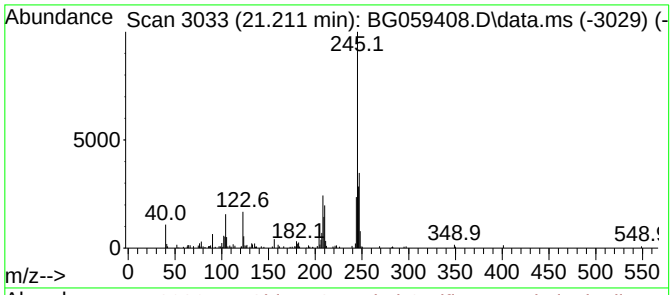
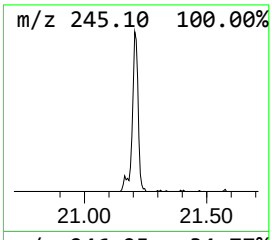
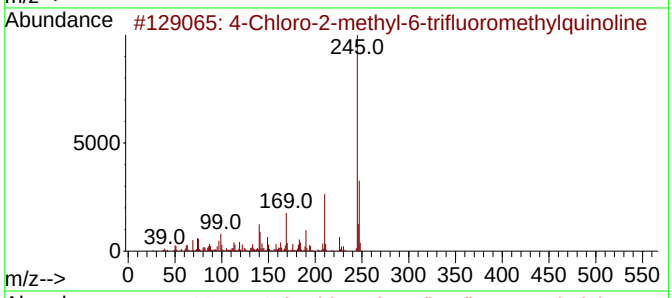
Peak Number 11 4-Chloro-2-methyl-6-trifluo... Concentration Rank 8

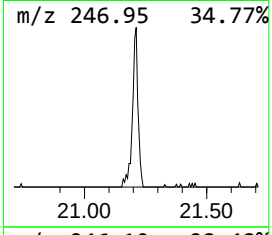
R.T.	EstConc	Area	Relative to ISTD	R.T.
21.211	5.13 ng/ul	177598	Chrysene-d12	21.910

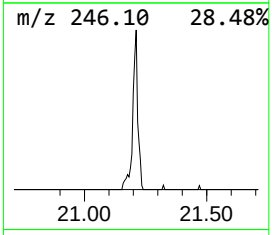
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Chloro-2-methyl-6-trifluoromet...	245	C11H7ClF3N	867167-05-3	53
2	2-(4-chlorophenyl)-5-fluoro-1H-i...	245	C14H9ClFN	881040-32-0	50
3	2-Chloro-4-methylpyridazino(6,1-...	245	C12H8ClN3O	001703-04-4	50
4	Quinoline, 2-(1H-benzimidazol-2-...	245	C16H11N3	014044-48-5	43
5	4-Chloro-2-methyl-7-(trifluorome...	245	C11H7ClF3N	018529-09-4	42

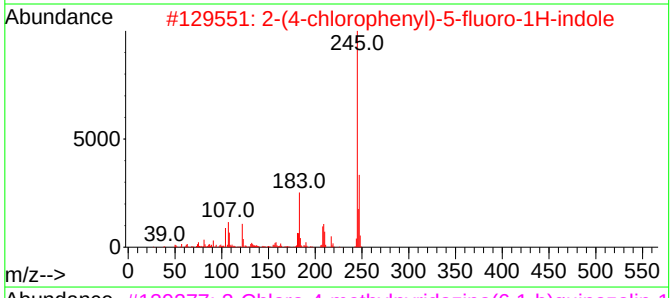


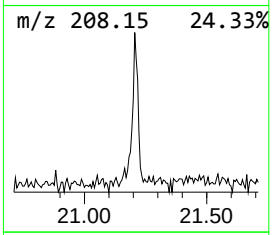


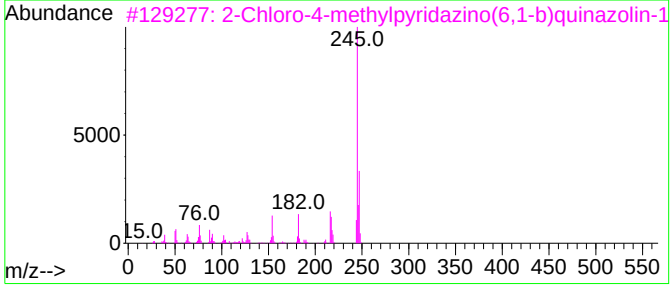


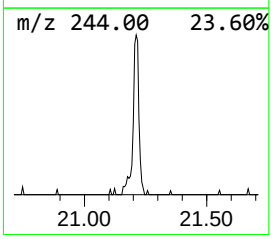












Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100923\
Data File : BG059408.D
Acq On : 9 Oct 2023 22:42
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Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown-01	4.836	208.9	ng/ul	2493980	1	8.226	238721	20.0
Butane, 2,3-dic...	5.136	17.4	ng/ul	208123	1	8.226	238721	20.0
unknown-02	5.359	9.5	ng/ul	113780	1	8.226	238721	20.0
unknown-03	7.633	3.5	ng/ul	41740	1	8.226	238721	20.0
unknown-04	8.297	3.3	ng/ul	39388	1	8.226	238721	20.0
Benzene,1-chlor...	9.501	23.4	ng/ul	279579	1	8.226	238721	20.0
unknown-05	10.847	5.3	ng/ul	101758	2	11.064	383001	20.0
Xenon	10.906	6.7	ng/ul	128479	2	11.064	383001	20.0
Nonanoic acid	11.787	2.0	ng/ul	39038	2	11.064	383001	20.0
unknown-06	17.845	6.8	ng/ul	238914	4	17.610	697114	20.0
4-Chloro-2-meth...	21.211	5.1	ng/ul	177598	5	21.910	693033	20.0