LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061423\
Data File : BG057705.D
Acq On : 14 Jun 2023 13:01
Operator : CG/JU
Sample : 03152-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RE137-CARBON-20230609

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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\8270-BG061323.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG057705.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.128	3	7	12	rVV	55995	72571	11.62%	1.012%
2	3.169	12	14	23	rVB3	10559	13612	2.18%	0.190%
3	3.357	41	46	62	rVB	194459	331242	53.06%	4.618%
4	3.486	64	68	76	rVB4	7544	12248	1.96%	0.171%
5	3.974	147	151	156	rBV	14958	23006	3.69%	0.321%
6	4.232	188	195	206	rBV	404235	624302	100.00%	8.703%
7	4.656	260	267	275	rBV2	12407	22586	3.62%	0.315%
8	4.896	301	308	321	rBV	341469	501955	80.40%	6.997%
9	5.043	328	333	340	rVB	30702	45525	7.29%	0.635%
10	5.331	378	382	391	rBV5	6978	13341	2.14%	0.186%
11	5.507	406	412	419	rBV	168175	255219	40.88%	3.558%
12	5.983	489	493	498	rVB3	5735	9601	1.54%	0.134%
13	6.095	506	512	518	rBV2	4510	8249	1.32%	0.115%
14	6.336	547	553	561	rBV	29699	46095	7.38%	0.643%
15	6.442	565	571	579	rVB	11714	18805	3.01%	0.262%
16	6.765	622	626	631	rVB2	5679	8222	1.32%	0.115%
17	7.035	666	672	675	rBV	8530	13485	2.16%	0.188%
18	7.088	675	681	689	rVB	175626	288334	46.19%	4.019%
19	7.317	714	720	726	rBV	4754	7868	1.26%	0.110%
20	7.940	816	826	832	rBV	132194	224366	35.94%	3.128%
21	8.093	848	852	857	rVB2	6565	9730	1.56%	0.136%
22	9.097	1017	1023	1031	rBV	91005	155081	24.84%	2.162%
23	9.332	1057	1063	1065	rBV2	3602	6672	1.07%	0.093%
24	9.379	1067	1071	1079	rVB	43362	71111	11.39%	0.991%
25	9.926	1159	1164	1168	rVB4	4121	6880	1.10%	0.096%
26	10.043	1180	1184	1188	rBV3	4140	6653	1.07%	0.093%
27	10.742	1293	1303	1310	rBV	177155	318291	50.98%	4.437%
28	10.854	1316	1322	1328	rVV	13826	22895	3.67%	0.319%
29	10.931	1330	1335	1341	rVB2	3631	6829	1.09%	0.095%
30	11.941	1500	1507	1513	rVB6	5224	10223	1.64%	0.143%
31	13.005	1680	1688	1693	rBV6	4987	9729	1.56%	0.136%
32	13.193	1714	1720	1727	rBV	238484	377789	60.51%	5.266%
33	13.281	1731	1735	1740	rVV4	3990	7127	1.14%	0.099%
34	13.345	1742	1746	1750	rVB5	5486	7861	1.26%	0.110%
35	13.486	1765	1770	1780	rVB	24611	38102	6.10%	0.531%
36	13.868	1831	1835	1841	rBV3	4344	7126	1.14%	0.099%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061423\
Data File : BG057705.D
Acq On : 14 Jun 2023 13:01
Operator : CG/JU
Sample : 03152-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RE137-CARBON-20230609

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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\8270-BG061323.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	14.573	1948	1955	1964	rVB	279589	444476	71.20%	6.196%
38	15.925	2177	2185	2189	rBV	58410	79100	12.67%	1.103%
39	16.054	2198	2207	2216	rBV	169323	264866	42.43%	3.692%
40	16.512	2280	2285	2289	rBV4	4886	7875	1.26%	0.110%
41	17.064	2375	2379	2385	rBV7	3270	6991	1.12%	0.097%
42	17.311	2415	2421	2429	rVB	392529	596353	95.52%	8.313%
43	18.204	2568	2573	2582	rBV2	44651	67678	10.84%	0.943%
44	18.281	2582	2586	2590	rVB	9430	13230	2.12%	0.184%
45	19.479	2786	2790	2796	rVB6	15456	18323	2.93%	0.255%
46	19.932	2861	2867	2872	rBV	473432	614075	98.36%	8.560%
47	20.243	2918	2920	2925	rVB3	5645	7173	1.15%	0.100%
48	20.743	3002	3005	3008	rVB2	7403	7595	1.22%	0.106%
49	21.230	3084	3088	3096	rBV3	73833	116497	18.66%	1.624%
50	21.307	3098	3101	3107	rVB	17695	22743	3.64%	0.317%
51	21.477	3126	3130	3134	rVB5	6623	9911	1.59%	0.138%
52	21.571	3140	3146	3155	rVB	371816	590397	94.57%	8.230%
53	21.783	3178	3182	3189	rVB3	10154	13247	2.12%	0.185%
54	22.388	3281	3285	3292	rVB6	7900	13326	2.13%	0.186%
55	22.746	3342	3346	3350	rVB5	7433	10729	1.72%	0.150%
56	23.069	3396	3401	3411	rVB6	8620	22078	3.54%	0.308%
57	23.862	3532	3536	3544	rVB6	5992	13946	2.23%	0.194%
58	24.732	3673	3684	3695	rBV	215848	609567	97.64%	8.497%
59	26.036	3903	3906	3917	rVB9	7509	20639	3.31%	0.288%

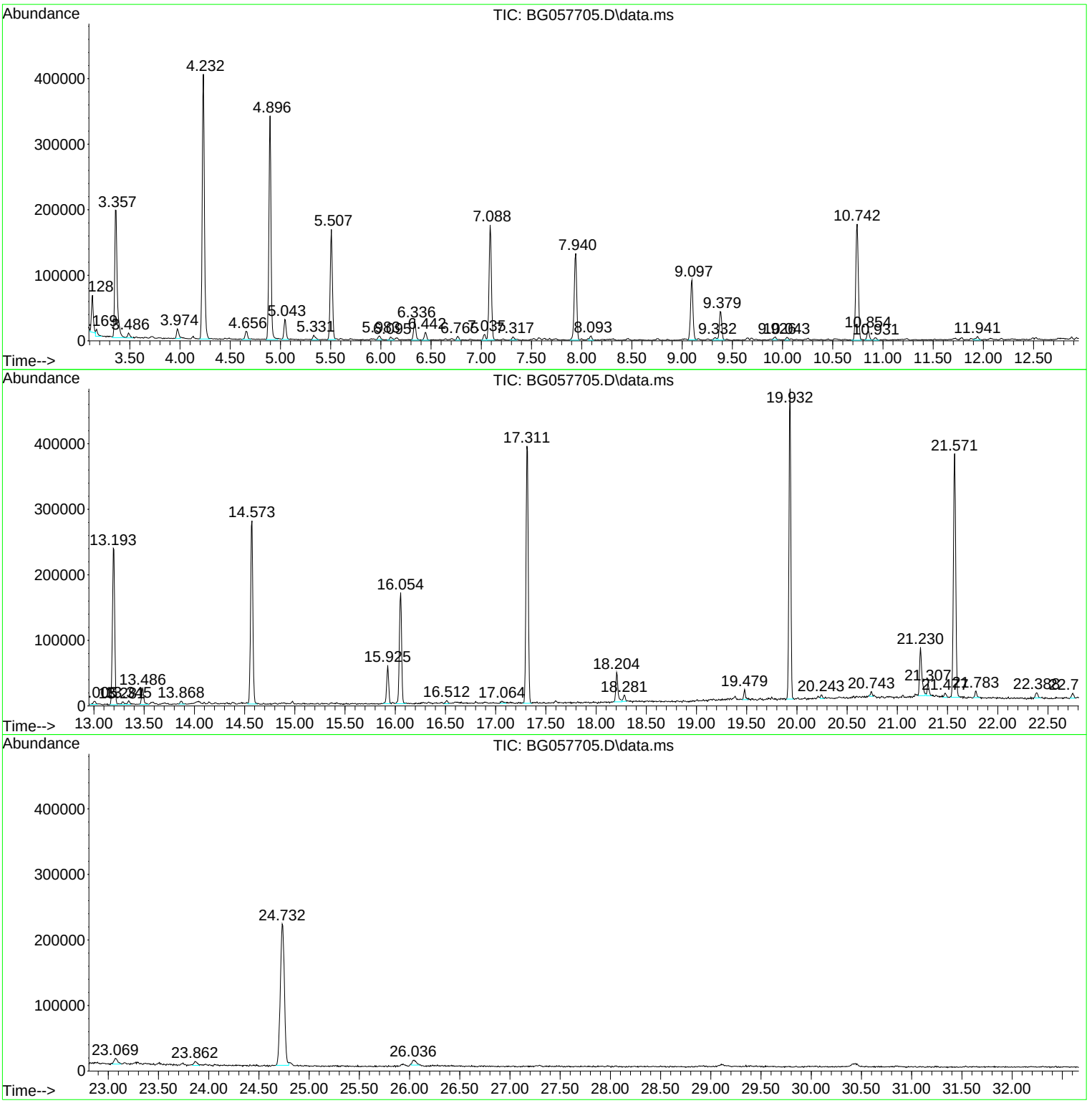
Sum of corrected areas: 7173546

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061423\
Data File : BG057705.D
Acq On : 14 Jun 2023 13:01
Operator : CG/JU
Sample : 03152-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RE137-CARBON-20230609

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061423\
Data File : BG057705.D
Acq On : 14 Jun 2023 13:01
Operator : CG/JU
Sample : 03152-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RE137-CARBON-20230609

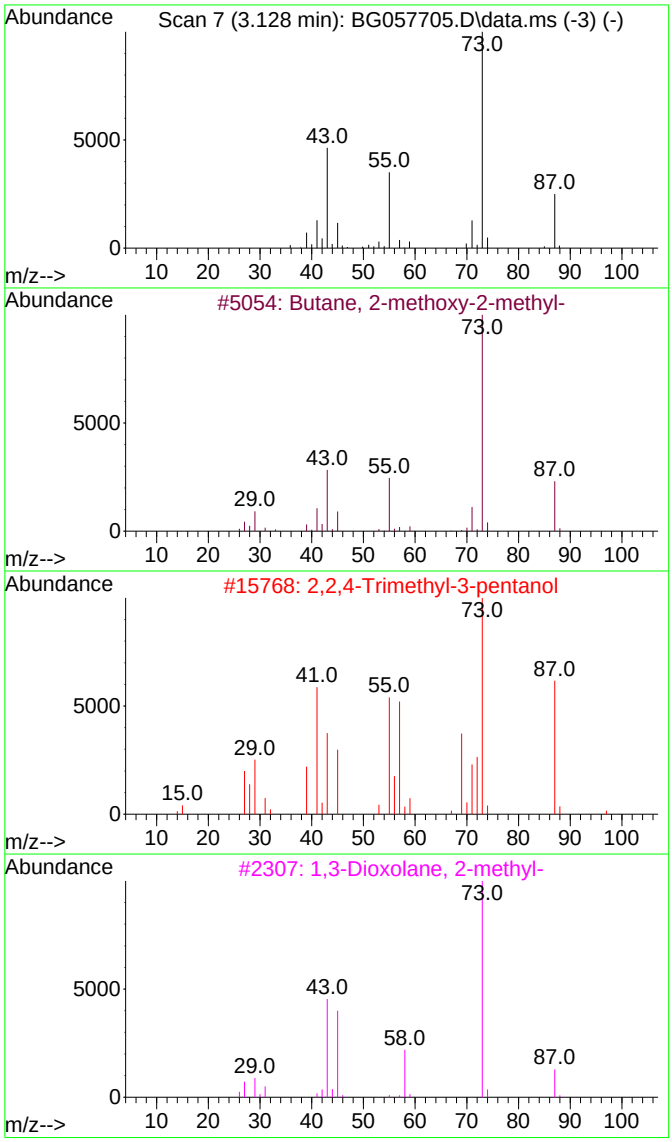
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

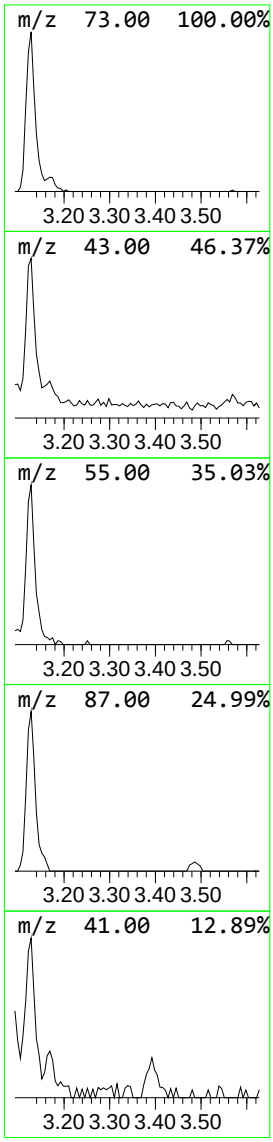
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.128	6.47 ng	72571	1,4-Dichlorobenzene-d4	7.940

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	38
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
5	Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061423\
Data File : BG057705.D
Acq On : 14 Jun 2023 13:01
Operator : CG/JU
Sample : 03152-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RE137-CARBON-20230609

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

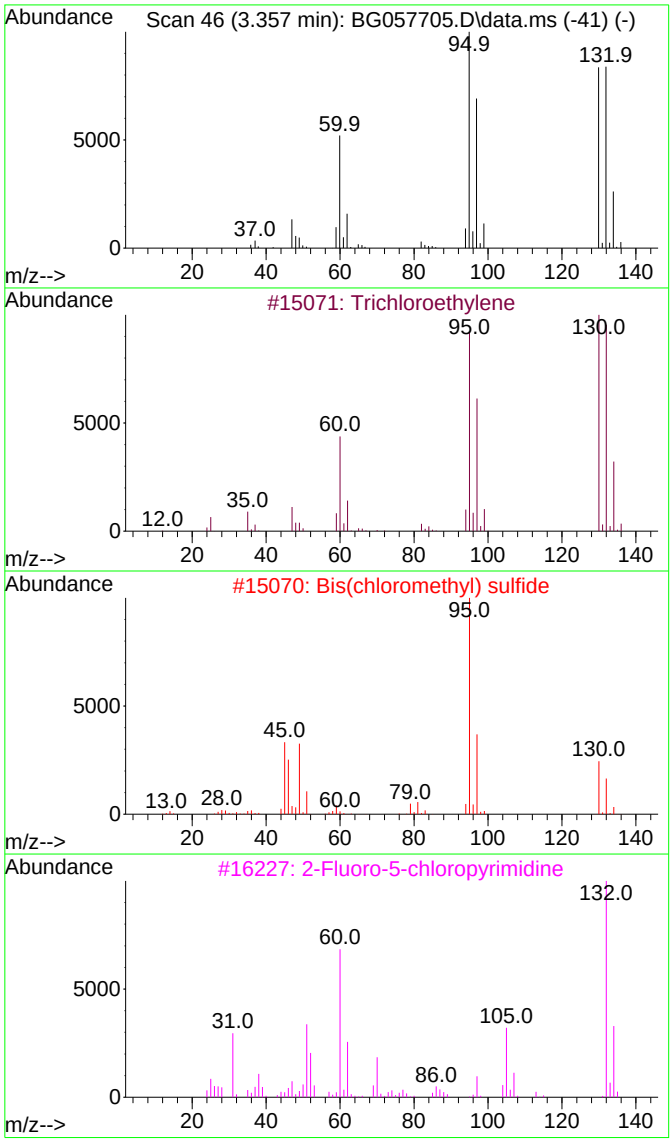
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

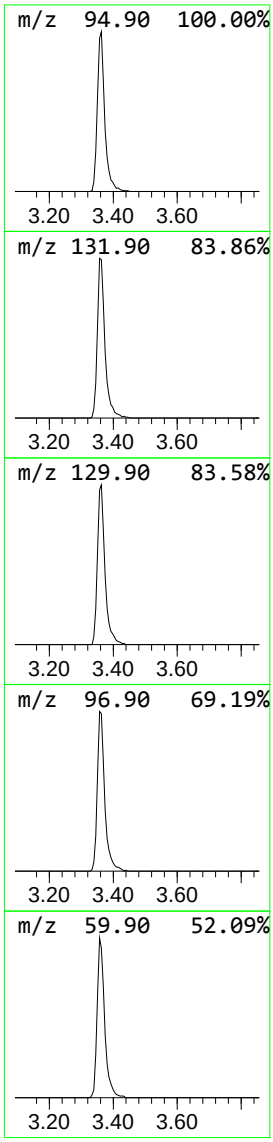
****

Peak Number 2 Trichloroethylene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.357	29.53 ng	331242	1,4-Dichlorobenzene-d4	7.940

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trichloroethylene	130	C2HCl3	000079-01-6	98
2	Bis(chloromethyl) sulfide	130	C2H4Cl2S	003592-44-7	49
3	2-Fluoro-5-chloropyrimidine	132	C4H2ClFN2	062802-37-3	25
4	Benzene, 1-chloro-4-fluoro-	130	C6H4ClF	000352-33-0	22
5	1H-Tetrazaborole, 5-chloro-4,5-d...	132	C2H6BClN4	021960-49-6	22





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061423\
Data File : BG057705.D
Acq On : 14 Jun 2023 13:01
Operator : CG/JU
Sample : 03152-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RE137-CARBON-20230609

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

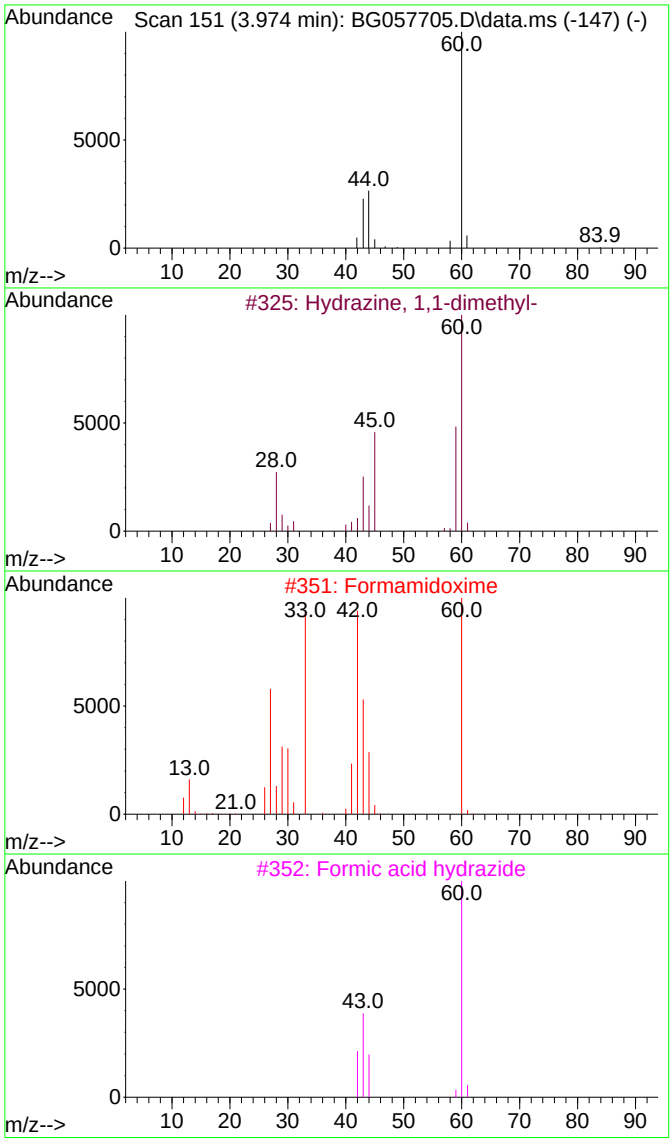
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

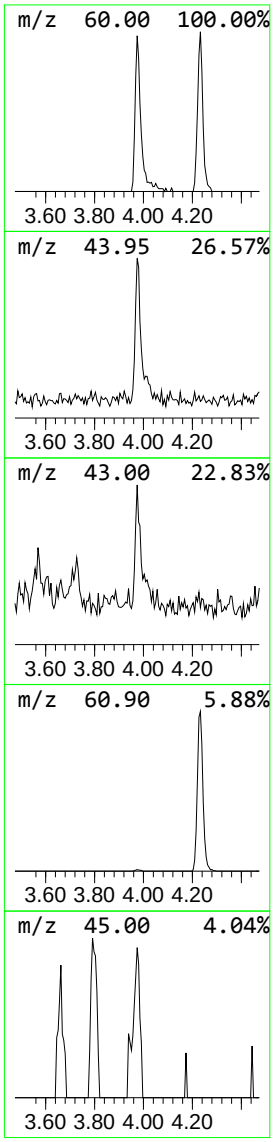
****

Peak Number 3 unknown3.974 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.974	2.05 ng	23006	1,4-Dichlorobenzene-d4	7.940

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydrazine, 1,1-dimethyl-	60	C2H8N2	000057-14-7	9
2	Formamidoxime	60	CH4N2O	000624-82-8	7
3	Formic acid hydrazide	60	CH4N2O	000624-84-0	7
4	Urea	60	CH4N2O	000057-13-6	5
5	Methyl formate	60	C2H4O2	000107-31-3	4





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061423\
Data File : BG057705.D
Acq On : 14 Jun 2023 13:01
Operator : CG/JU
Sample : 03152-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RE137-CARBON-20230609

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

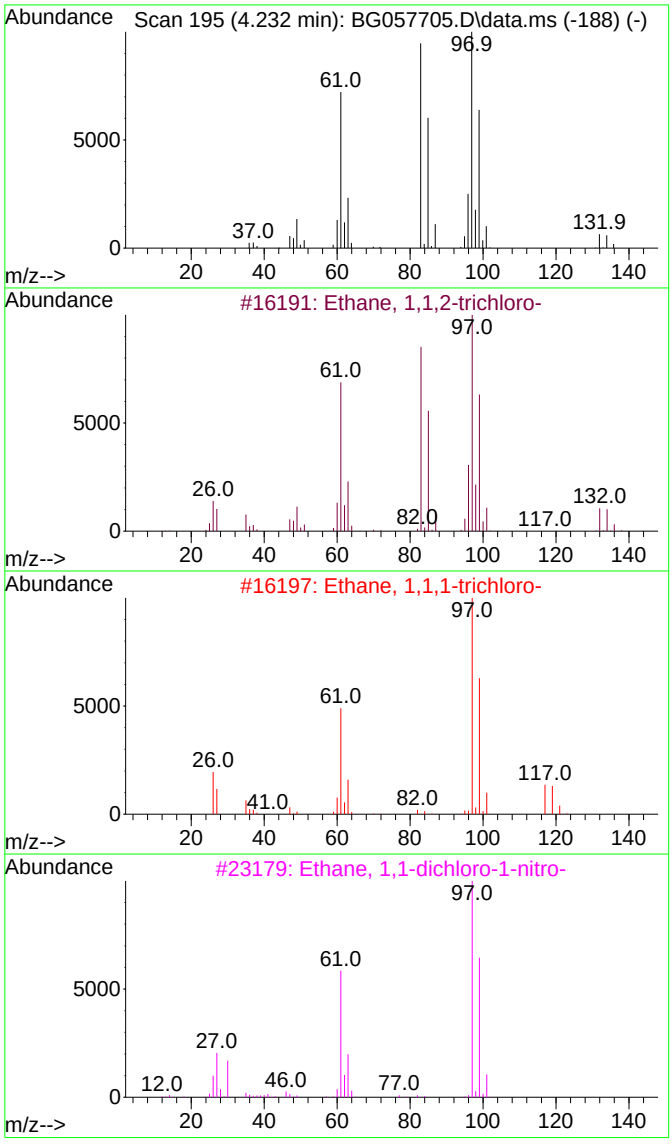
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

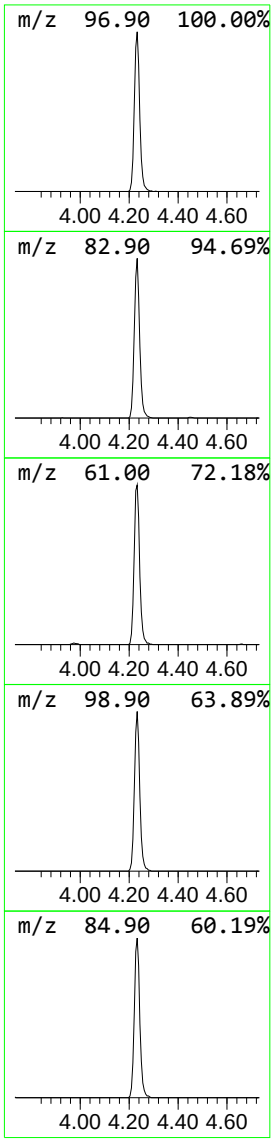
****

Peak Number 4 Ethane, 1,1,2-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.232	55.65 ng	624302	1,4-Dichlorobenzene-d4	7.940

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethane, 1,1,2-trichloro-	132	C2H3Cl3	000079-00-5	98
2	Ethane, 1,1,1-trichloro-	132	C2H3Cl3	000071-55-6	43
3	Ethane, 1,1-dichloro-1-nitro-	143	C2H3Cl2NO2	000594-72-9	43
4	Propanoyl chloride, 2,2-dichloro-	160	C3H3Cl3O	026073-26-7	32
5	Propane, 1,2,2-trichloro-	146	C3H5Cl3	003175-23-3	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061423\
Data File : BG057705.D
Acq On : 14 Jun 2023 13:01
Operator : CG/JU
Sample : 03152-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RE137-CARBON-20230609

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

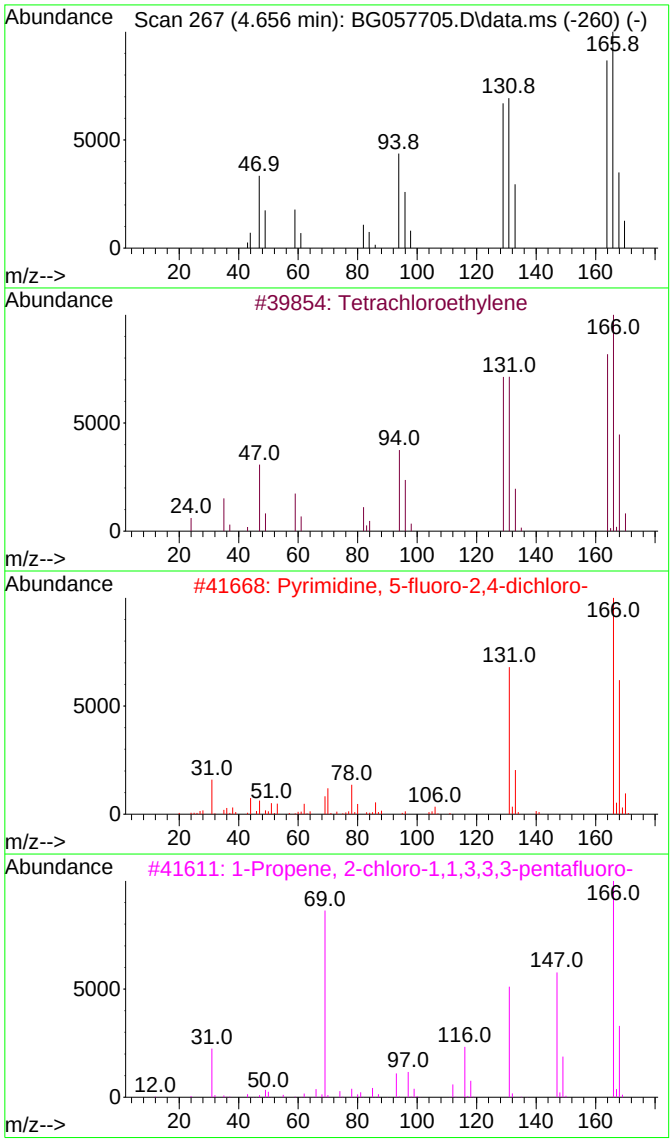
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

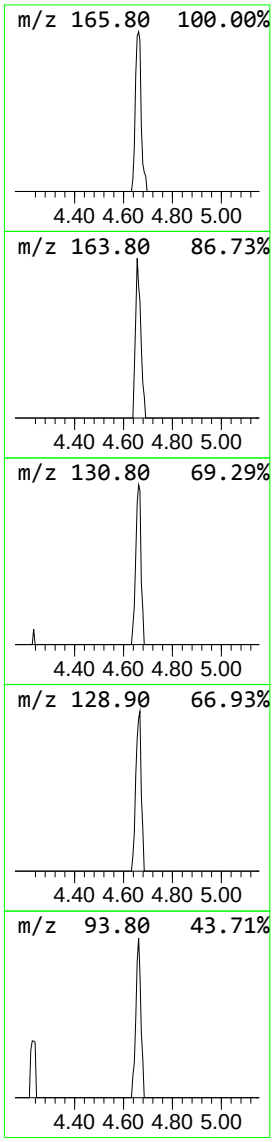
****

Peak Number 5 Tetrachloroethylene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.656	2.01 ng	22586	1,4-Dichlorobenzene-d4	7.940

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetrachloroethylene	164	C2Cl4	000127-18-4	96
2	Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HCl2FN2	002927-71-1	16
3	1-Propene, 2-chloro-1,1,3,3,3-pe...	166	C3ClF5	002804-50-4	9
4	4,5-Dichlorocyclopent-4-ene-1,3-...	164	C5H2Cl2O2	003229-28-5	9
5	4,6-Dichloro-2-fluoropyrimidine	166	C4HCl2FN2	003824-45-1	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061423\
Data File : BG057705.D
Acq On : 14 Jun 2023 13:01
Operator : CG/JU
Sample : 03152-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RE137-CARBON-20230609

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

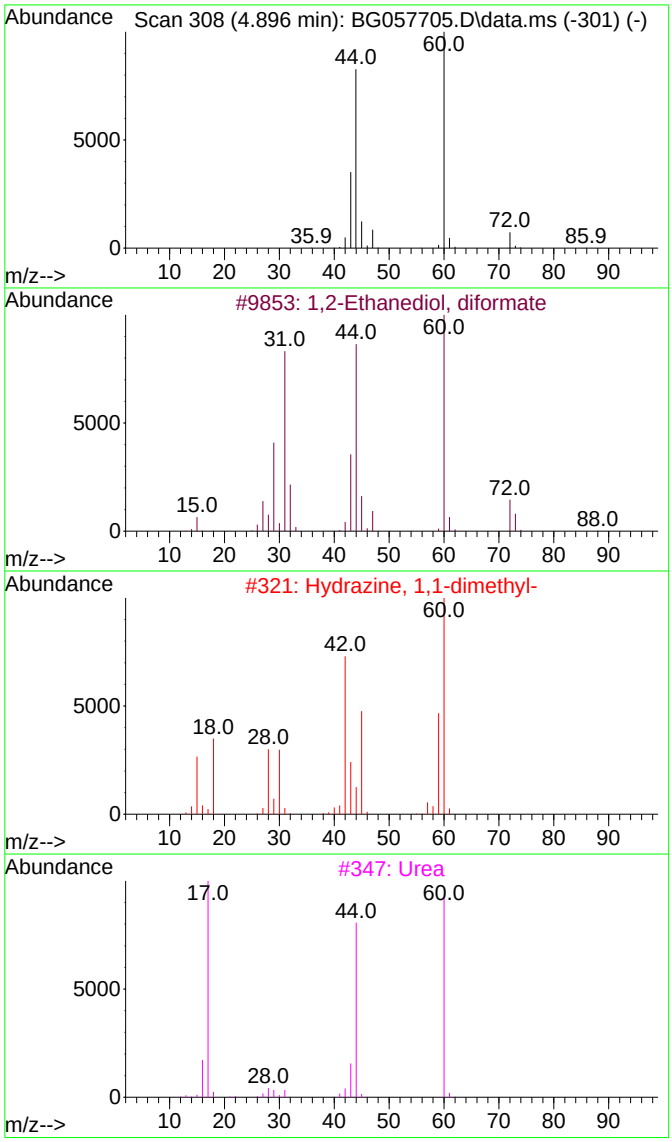
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

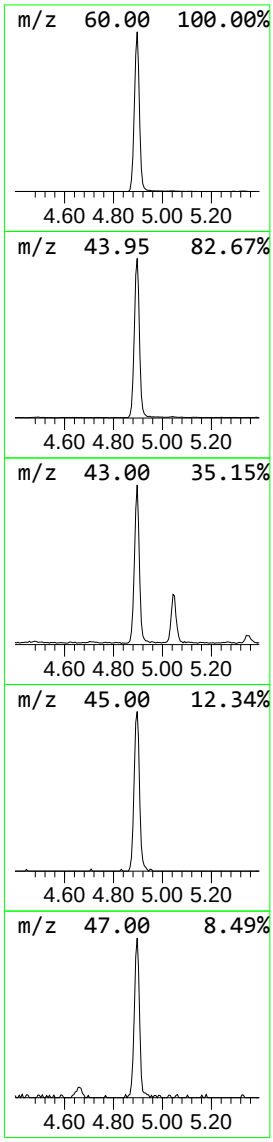
****

Peak Number 6 unknown4.896 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.896	44.74 ng	501955	1,4-Dichlorobenzene-d4	7.940

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Ethanediol, diformate	118	C4H6O4	000629-15-2	37
2	Hydrazine, 1,1-dimethyl-	60	C2H8N2	000057-14-7	9
3	Urea	60	CH4N2O	000057-13-6	9
4	Propanedioic acid, propyl-	146	C6H10O4	000616-62-6	9
5	Acetic acid, hydroxy[(1-oxo-2-pr...	145	C5H7NO4	006737-24-2	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061423\
Data File : BG057705.D
Acq On : 14 Jun 2023 13:01
Operator : CG/JU
Sample : 03152-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RE137-CARBON-20230609

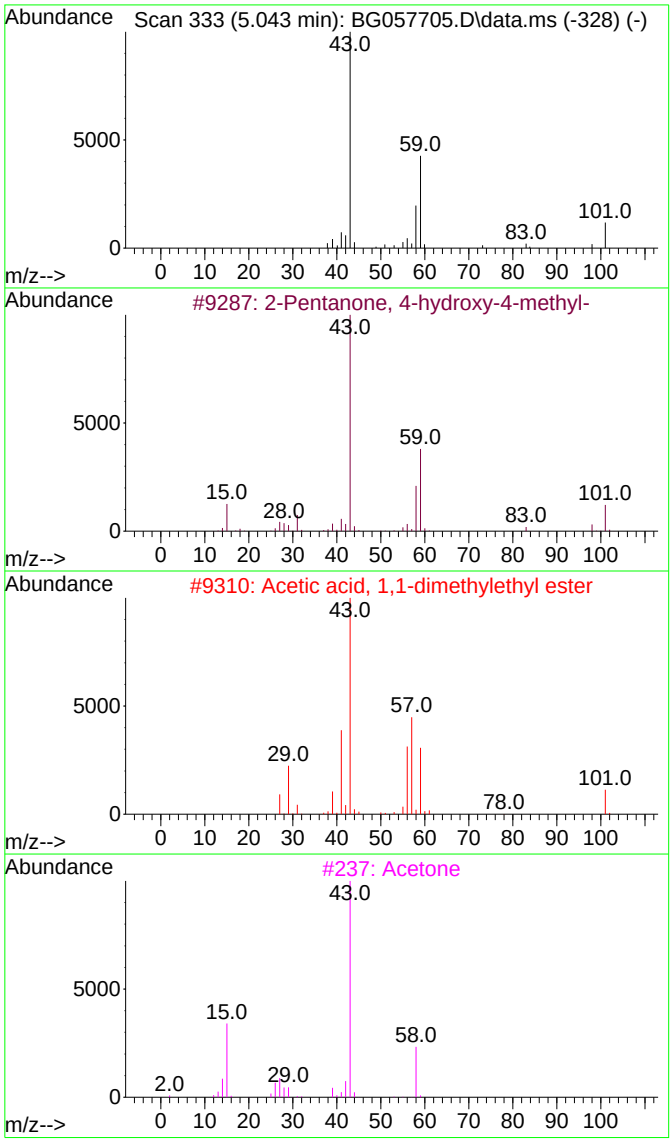
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

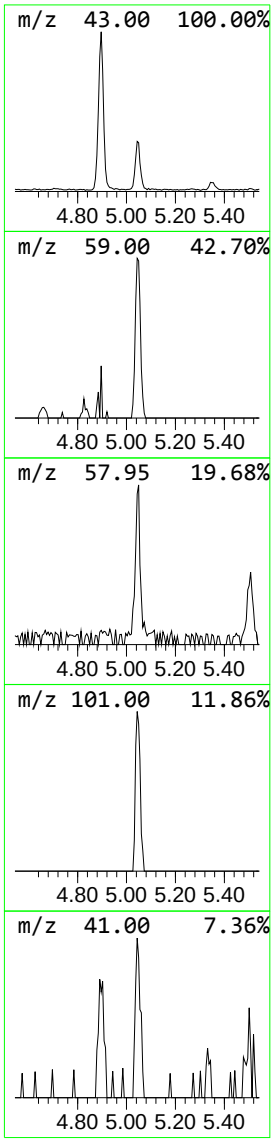
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.043	4.06 ng	45525	1,4-Dichlorobenzene-d4	7.940

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	36
3	Acetone	58	C3H6O	000067-64-1	16
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	14
5	N-(n-Propyl)acetamide	101	C5H11NO	005331-48-6	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061423\
Data File : BG057705.D
Acq On : 14 Jun 2023 13:01
Operator : CG/JU
Sample : 03152-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RE137-CARBON-20230609

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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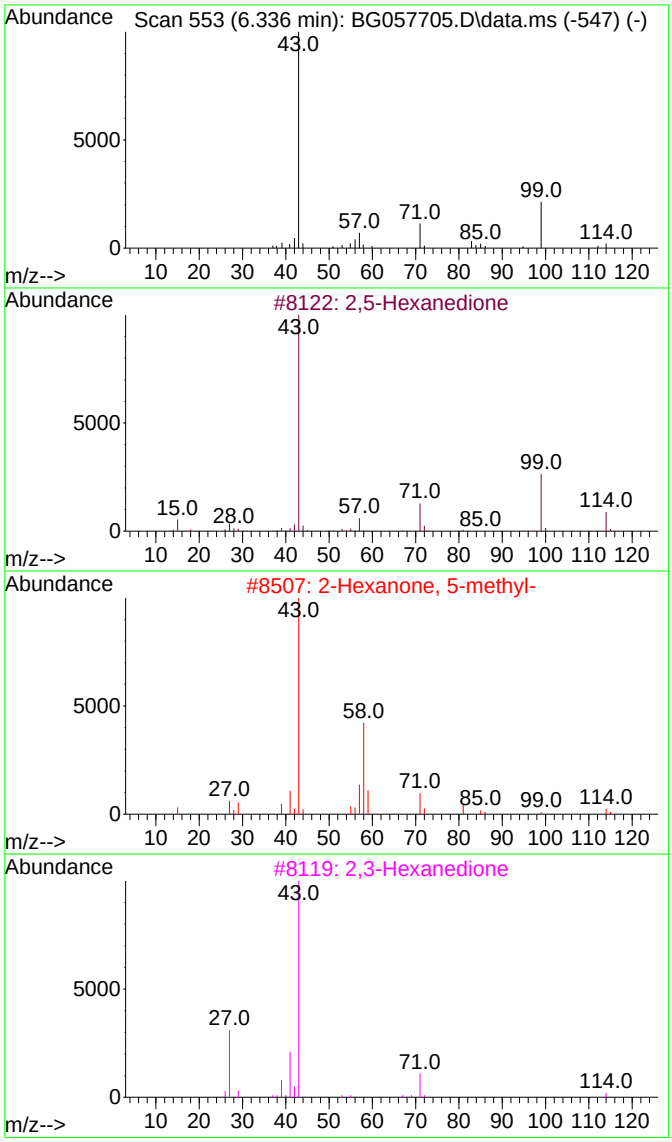
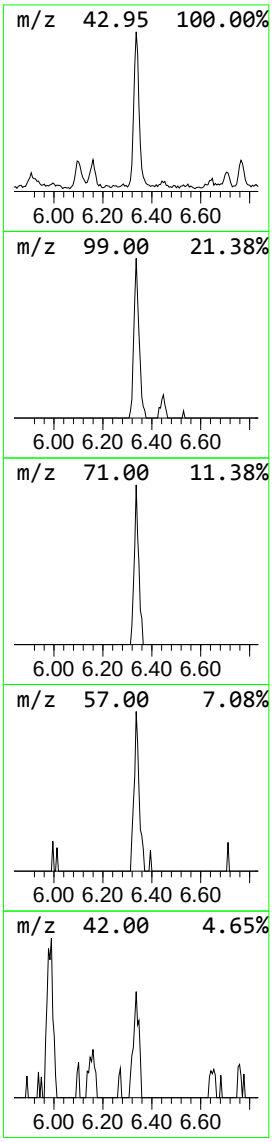
Peak Number 8 2,5-Hexanedione Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.336	4.11 ng	46095	1,4-Dichlorobenzene-d4	7.940

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,5-Hexanedione	114	C6H10O2	000110-13-4	64
2	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	37
3	2,3-Hexanedione	114	C6H10O2	003848-24-6	35
4	2-Hexanone, 4-methyl-	114	C7H14O	000105-42-0	25
5	2-Pentanone, 3-ethyl-	114	C7H14O	006137-03-7	23





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061423\
Data File : BG057705.D
Acq On : 14 Jun 2023 13:01
Operator : CG/JU
Sample : 03152-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RE137-CARBON-20230609

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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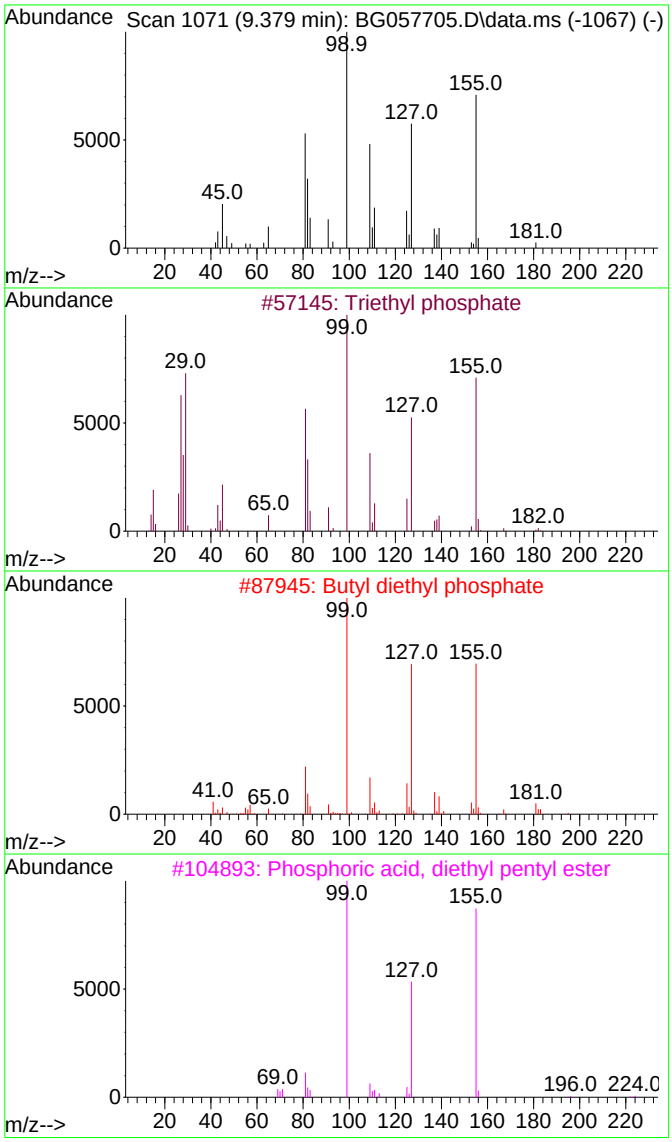
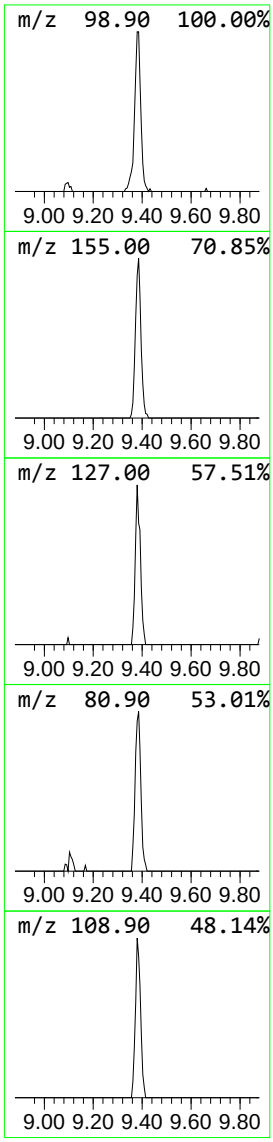
Peak Number 9 Triethyl phosphate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.379	4.47 ng	71111	Naphthalene-d8	10.743

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triethyl phosphate	182	C6H15O4P	000078-40-0	80
2	Butyl diethyl phosphate	210	C8H19O4P	1000298-58-0	64
3	Phosphoric acid, diethyl pentyl ...	224	C9H21O4P	020195-08-8	36
4	2H-Pyran-2-carboxylic acid, 6-et...	200	C10H16O4	013687-98-4	22
5	Benzene, 2-fluoro-1-methyl-4-nitro-	155	C7H6FNO2	001427-07-2	12





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Data File : BG057705.D
Acq On : 14 Jun 2023 13:01
Operator : CG/JU
Sample : 03152-01
Misc :
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Instrument :
BNA_G
ClientSampleId :
RE137-CARBON-20230609

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M
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TIC Library : C:\Database\NIST20.L
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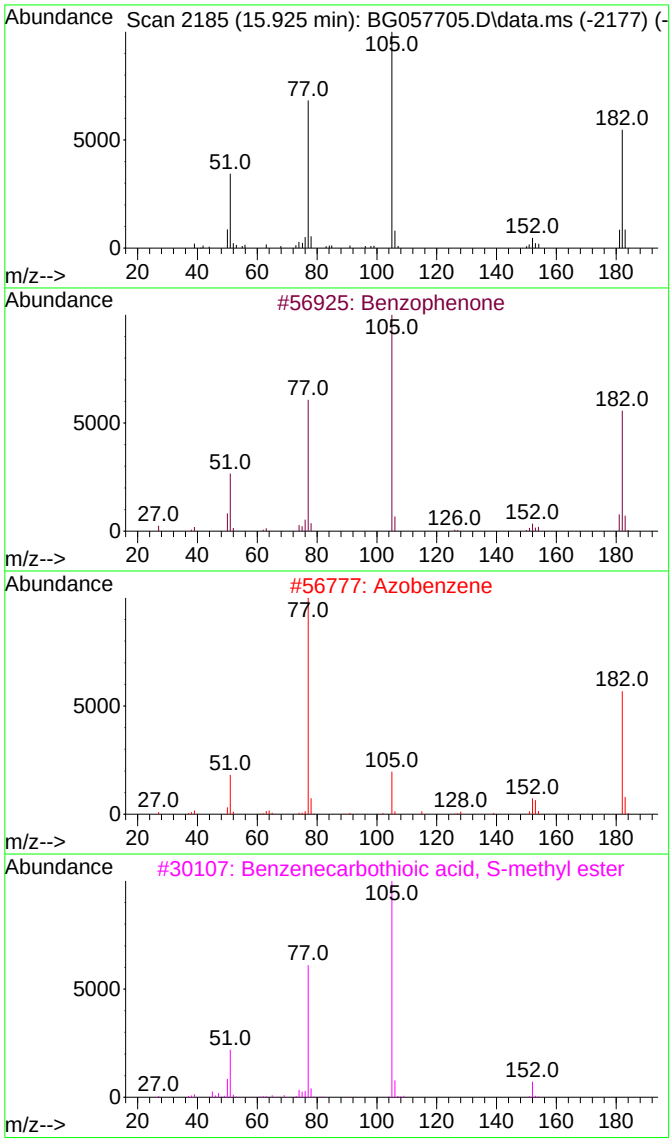
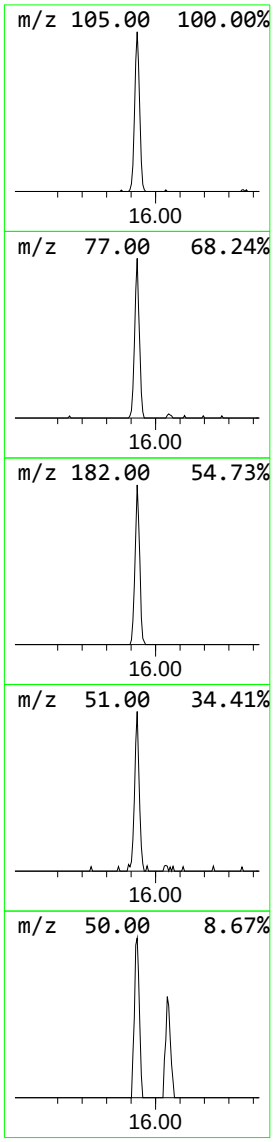
Peak Number 10 Benzophenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.925	3.56 ng	79100	Acenaphthene-d10	14.573

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	Azobenzene	182	C12H10N2	000103-33-3	64
3	Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	47
4	1-Propanone, 2-bromo-1-phenyl-	212	C9H9BrO	002114-00-3	47
5	2-Phenyl-4-ethylidene-2-oxazolin...	187	C11H9NO2	037791-41-6	47





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Acq On : 14 Jun 2023 13:01
Operator : CG/JU
Sample : 03152-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M
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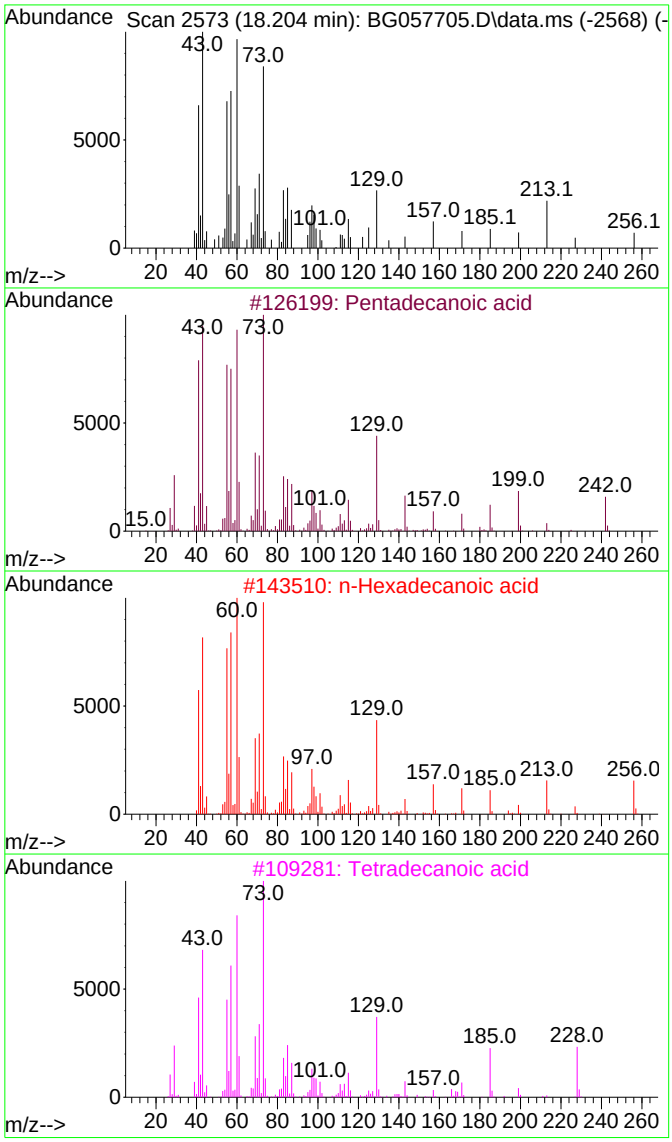
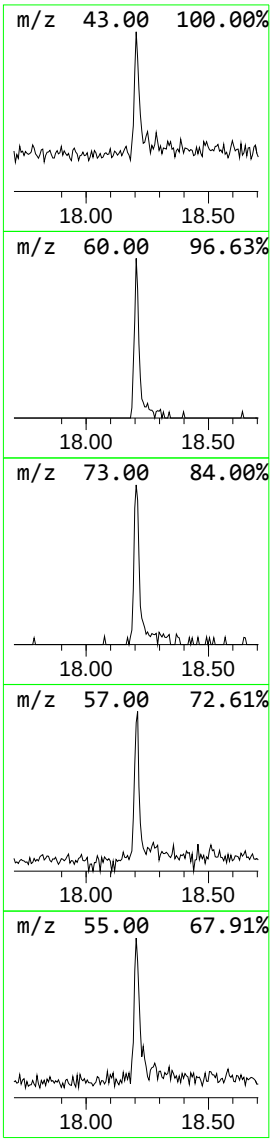
Peak Number 11 Pentadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.204	2.27 ng	67678	Phenanthrene-d10	17.311

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	53
4	Dodecanoic acid	200	C12H24O2	000143-07-7	53
5	Tridecanoic acid	214	C13H26O2	000638-53-9	50





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Operator : CG/JU
Sample : 03152-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
RE137-CARBON-20230609

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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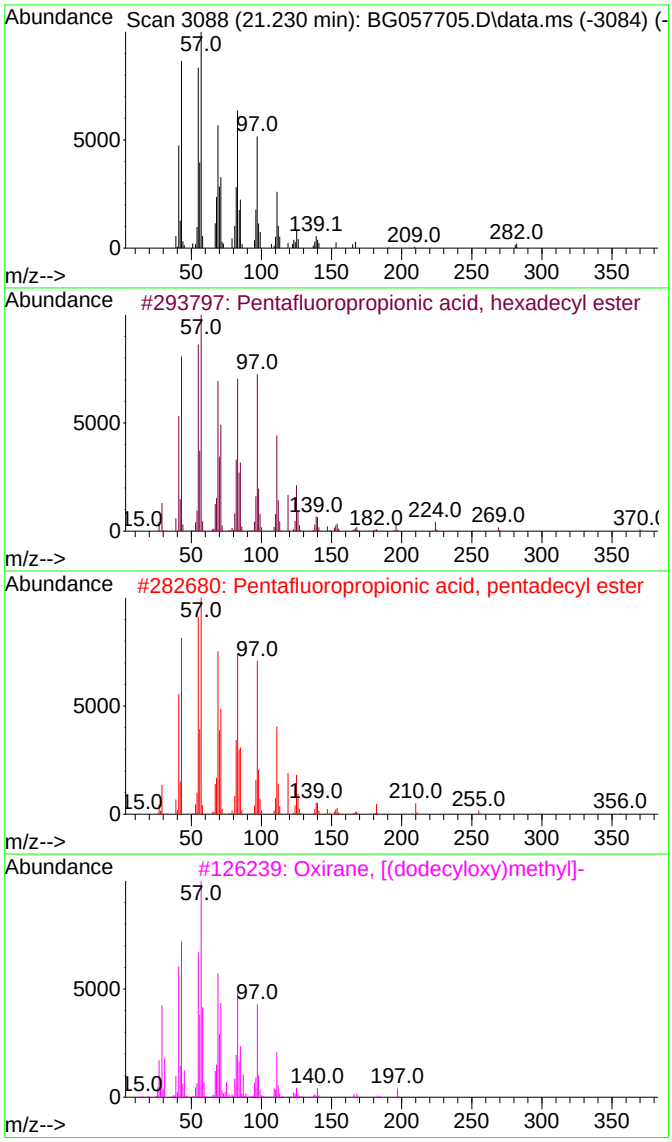
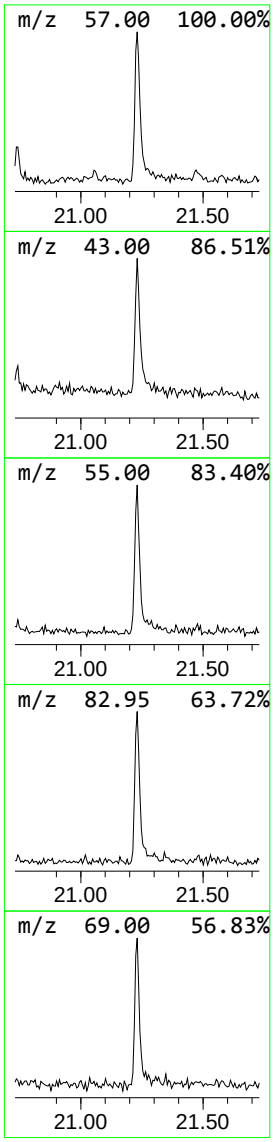
Peak Number 12 Pentafluoropropionic acid, ... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.230	3.95 ng	116497	Chrysene-d12	21.571

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	91
2	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	91
3	Oxirane, [(dodecyloxy)methyl]-	242	C15H30O2	002461-18-9	91
4	1-Octadecanol	270	C18H38O	000112-92-5	91
5	Heptacosyl acetate	438	C29H58O2	1000351-78-2	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061423\
Data File : BG057705.D
Acq On : 14 Jun 2023 13:01
Operator : CG/JU
Sample : 03152-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
Butane, 2-metho...	3.128	6.5	ng	72571	1	7.940	224366	20.0
Trichloroethylene	3.357	29.5	ng	331242	1	7.940	224366	20.0
unknown3.974	3.974	2.0	ng	23006	1	7.940	224366	20.0
Ethane, 1,1,2-t...	4.232	55.6	ng	624302	1	7.940	224366	20.0
Tetrachloroethy...	4.656	2.0	ng	22586	1	7.940	224366	20.0
unknown4.896	4.896	44.7	ng	501955	1	7.940	224366	20.0
2-Pentanone, 4-...	5.043	4.1	ng	45525	1	7.940	224366	20.0
2,5-Hexanedione	6.336	4.1	ng	46095	1	7.940	224366	20.0
Triethyl phosphate	9.379	4.5	ng	71111	2	10.743	318291	20.0
Benzophenone	15.925	3.6	ng	79100	3	14.573	444476	20.0
Pentadecanoic acid	18.204	2.3	ng	67678	4	17.311	596353	20.0
Pentafluoroprop...	21.230	4.0	ng	116497	5	21.571	590397	20.0