LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012721\
Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : 0051-045-COMP-E-ELUTRIATE

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-BG012521.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG048042.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.678	390	398	407	rBV	559203	882581	22.42%	4.231%
2	7.270	660	669	680	rBV	380680	718227	18.24%	3.443%
3	8.122	808	814	821	rBV	173085	290754	7.39%	1.394%
4	9.286	1004	1012	1023	rVB	561290	1005718	25.55%	4.821%
5	10.937	1283	1293	1302	rBV	218228	400017	10.16%	1.918%
6	13.375	1700	1708	1714	rBV	1537302	2299101	58.40%	11.022%
7	14.192	1841	1847	1851	rBV2	42712	66248	1.68%	0.318%
8	14.374	1874	1878	1885	rBV5	26241	44830	1.14%	0.215%
9	14.744	1935	1941	1951	rVB	379055	593689	15.08%	2.846%
10	15.543	2071	2077	2082	rBV3	22119	43459	1.10%	0.208%
11	16.230	2185	2194	2200	rBV	1368762	2102232	53.40%	10.078%
12	17.394	2385	2392	2396	rBV3	28155	40277	1.02%	0.193%
13	17.488	2401	2408	2414	rVV	506320	769731	19.55%	3.690%
14	17.764	2447	2455	2466	rBV	963021	1253907	31.85%	6.011%
15	18.369	2553	2558	2566	rBV	535481	770146	19.56%	3.692%
16	18.440	2566	2570	2576	rVB	78653	111634	2.84%	0.535%
17	18.880	2641	2645	2650	rBV	35065	56543	1.44%	0.271%
18	19.139	2684	2689	2694	rBV	320626	371672	9.44%	1.782%
19	19.221	2699	2703	2707	rBV	97329	120853	3.07%	0.579%
20	19.292	2711	2715	2721	rBV2	33662	44656	1.13%	0.214%
21	19.632	2768	2773	2784	rBV	430264	623316	15.83%	2.988%
22	19.767	2792	2796	2803	rVB	101437	129408	3.29%	0.620%
23	20.091	2845	2851	2861	rVB	2971792	3936934	100.00%	18.873%
24	20.296	2882	2886	2890	rBV	73383	86731	2.20%	0.416%
25	20.772	2964	2967	2974	rVB	150815	203843	5.18%	0.977%
26	20.843	2974	2979	2983	rBV4	31131	65418	1.66%	0.314%
27	20.990	2999	3004	3009	rBV	670737	829468	21.07%	3.976%
28	21.042	3009	3013	3017	rVV	160911	209275	5.32%	1.003%
29	21.107	3021	3024	3035	rVB	60277	99303	2.52%	0.476%
30	21.395	3069	3073	3081	rBV	310263	457835	11.63%	2.195%
31	21.554	3096	3100	3104	rVB	75979	101949	2.59%	0.489%
32	21.624	3110	3112	3121	rVB4	34218	76706	1.95%	0.368%
33	21.759	3129	3135	3145	rVB2	558802	987511	25.08%	4.734%
34	23.040	3350	3353	3359	rVB5	29265	46244	1.17%	0.222%
35	25.050	3685	3695	3704	rBV2	316074	904730	22.98%	4.337%
36	26.383	3918	3922	3935	rVB8	38871	114824	2.92%	0.550%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012721\
Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
0051-045-COMP-E-ELUTRIATE

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-BG012521.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

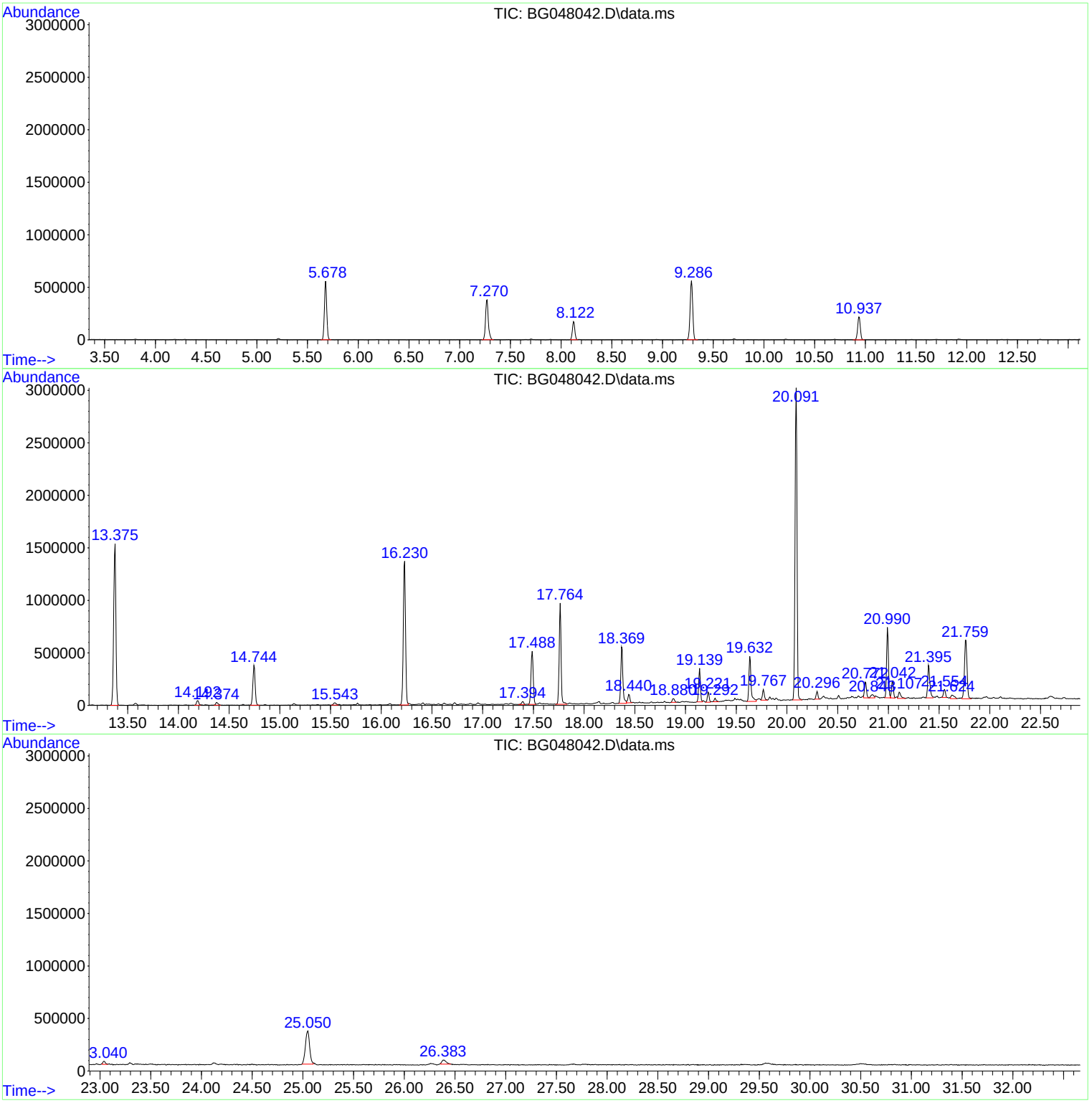
Sum of corrected areas: 20859770

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012721\
Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
0051-045-COMP-E-ELUTRIATE

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012721\
Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
0051-045-COMP-E-ELUTRIATE

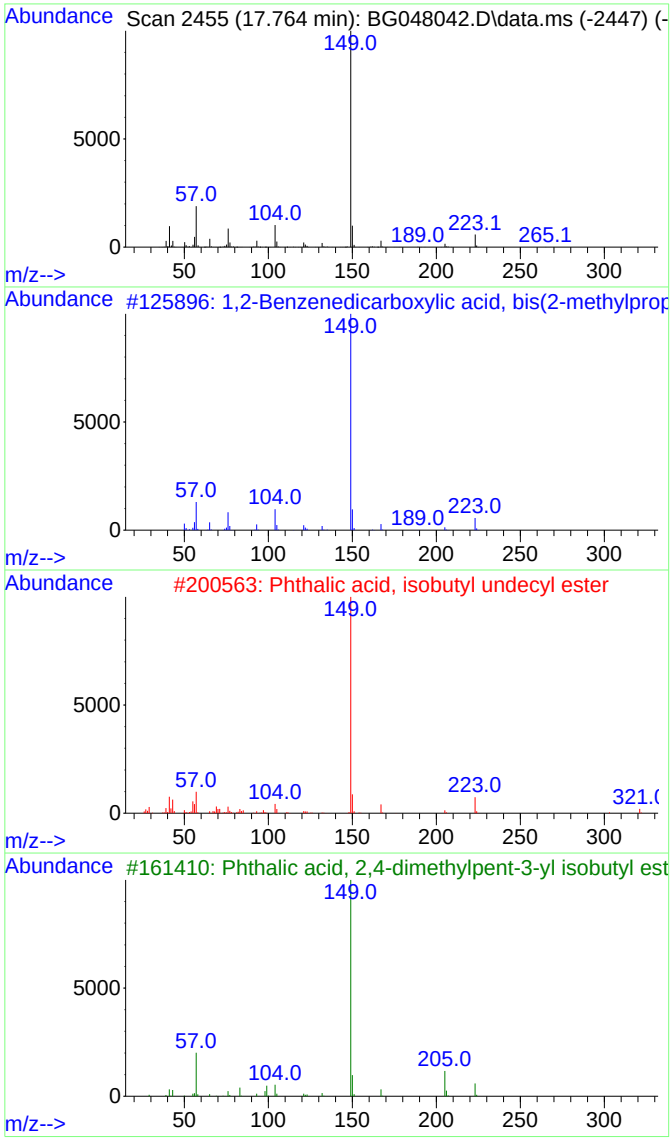
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

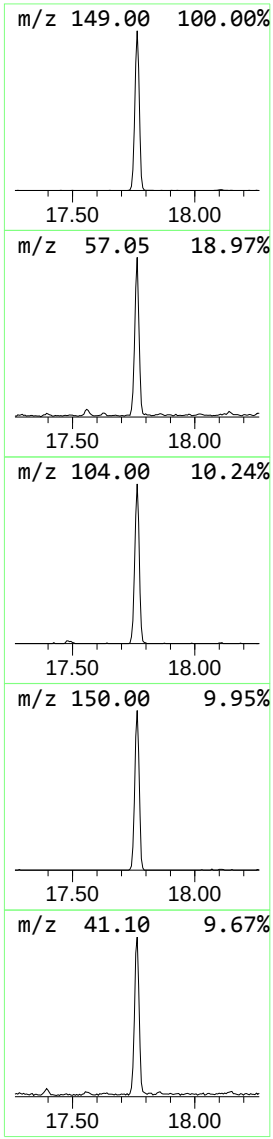
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 1,2-Benzenedicarboxylic aci... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.764	32.58 ng	1253910	Phenanthrene-d10	17.488

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Benzenedicarboxylic acid, bi...	278	C16H22O4	000084-69-5	91
2	Phthalic acid, isobutyl undecyl ...	376	C23H36O4	1000308-97-3	83
3	Phthalic acid, 2,4-dimethylpent-...	320	C19H28O4	1000356-84-3	83
4	Phthalic acid, 6-ethyl-3-octyl i...	362	C22H34O4	1000315-17-9	83
5	Phthalic acid, isobutyl octyl ester	334	C20H30O4	1000309-04-5	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012721\
Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
0051-045-COMP-E-ELUTRIATE

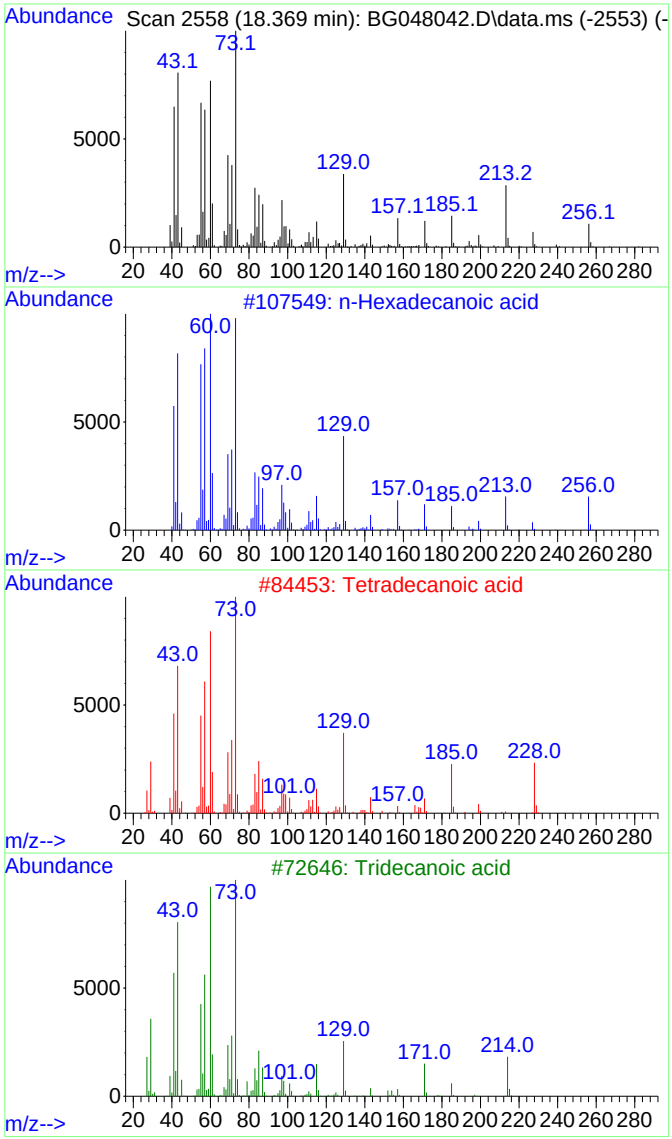
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

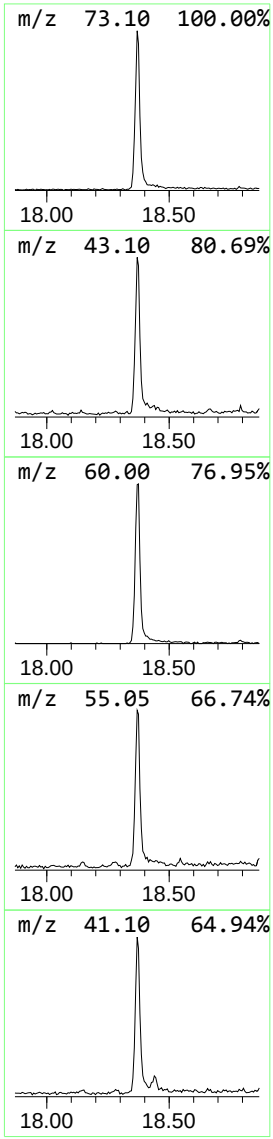
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.369	20.01 ng	770146	Phenanthrene-d10	17.488

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	93
3	Tridecanoic acid	214	C13H26O2	000638-53-9	90
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5	n-Decanoic acid	172	C10H20O2	000334-48-5	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012721\
Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
0051-045-COMP-E-ELUTRIATE

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

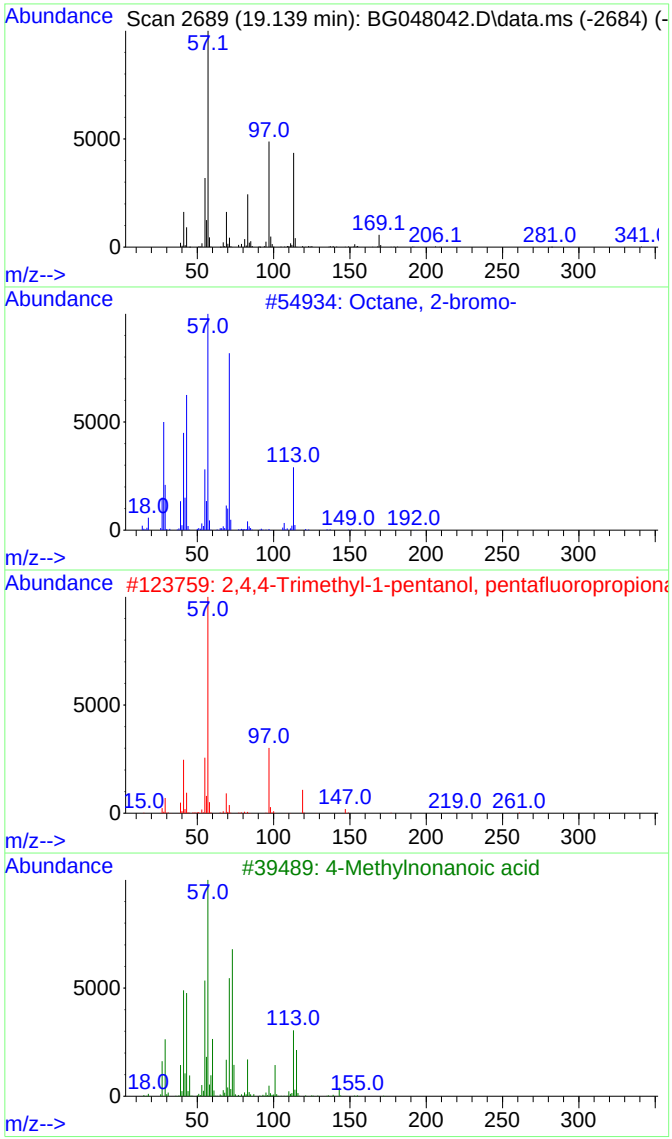
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

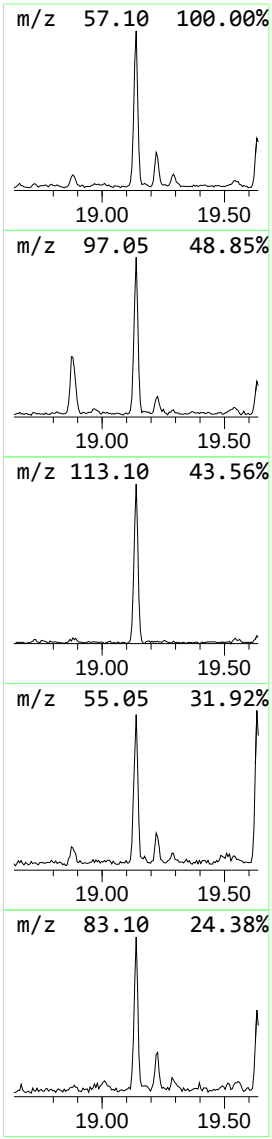
****

Peak Number 3 unknown19.139 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.139	9.66 ng	371672	Phenanthrene-d10	17.488

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2-bromo-	192	C8H17Br	000557-35-7	38
2	2,4,4-Trimethyl-1-pentanol, pent...	276	C11H17F5O2	1000365-51-0	38
3	4-Methylnonanoic acid	172	C10H20O2	045019-28-1	28
4	1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	28
5	2,4,4-Trimethyl-1-pentanol, hept...	326	C12H17F7O2	1000365-01-4	27





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012721\
Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
0051-045-COMP-E-ELUTRIATE

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

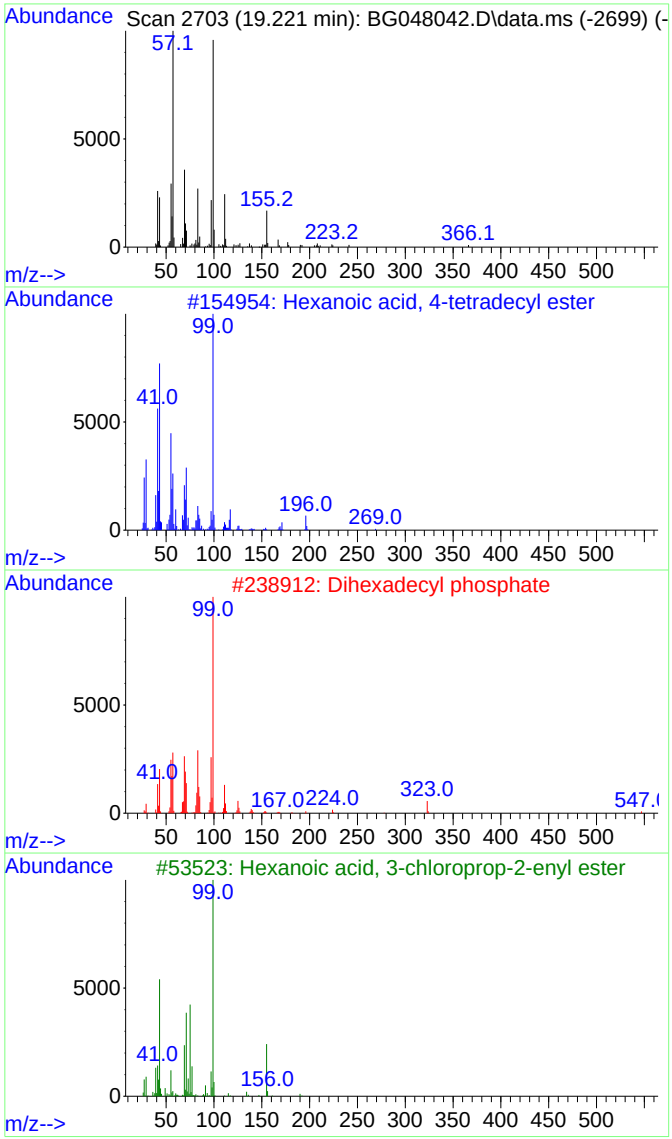
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

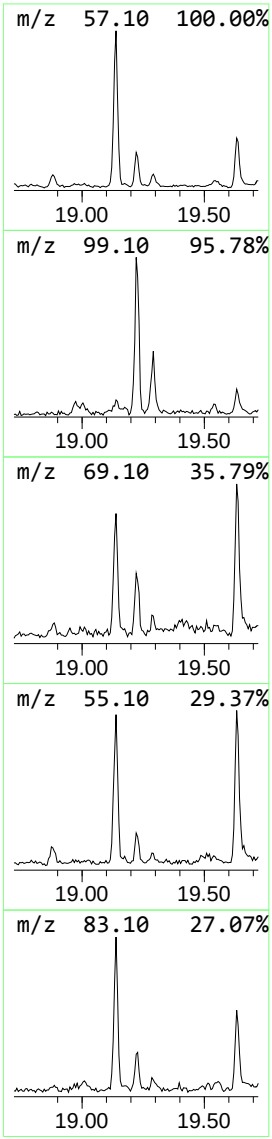
****

Peak Number 4 unknown19.221 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.221	3.14 ng	120853	Phenanthrene-d10	17.488

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 4-tetradecyl ester	312	C20H40O2	1000279-29-8	43
2	Dihexadecyl phosphate	546	C32H67O4P	002197-63-9	40
3	Hexanoic acid, 3-chloroprop-2-en...	190	C9H15ClO2	1000299-36-1	38
4	2(3H)-Furanone, 5-butyldihydro-4...	156	C9H16O2	055013-32-6	38
5	trans-3-Methyl-4-octanolide	156	C9H16O2	039638-67-0	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012721\
Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
0051-045-COMP-E-ELUTRIATE

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

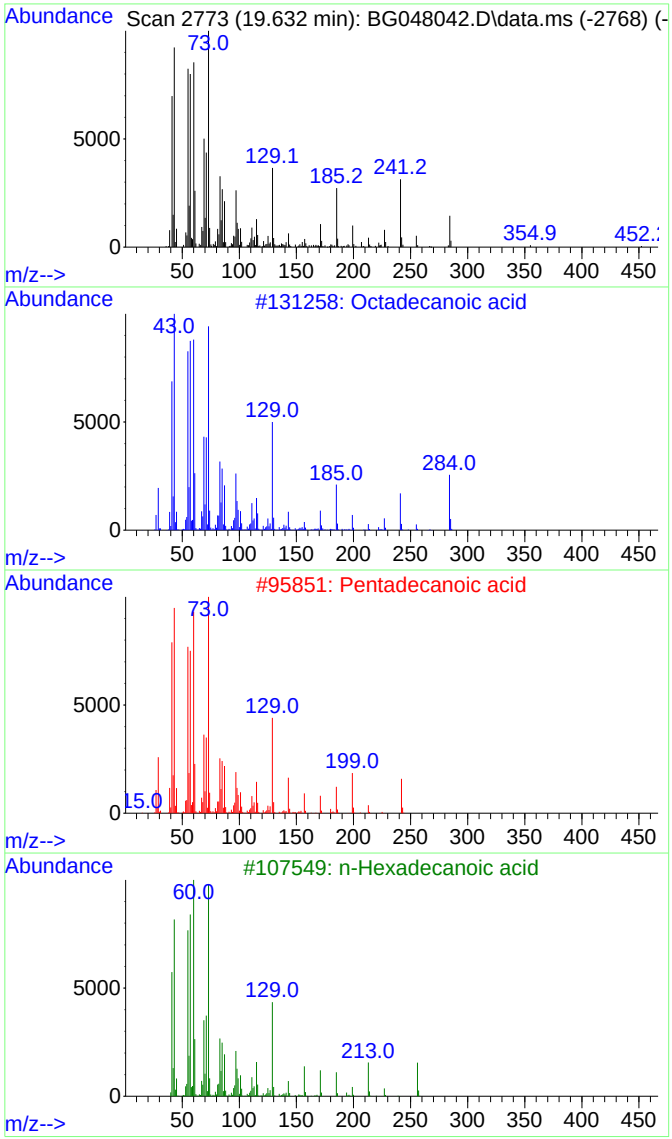
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

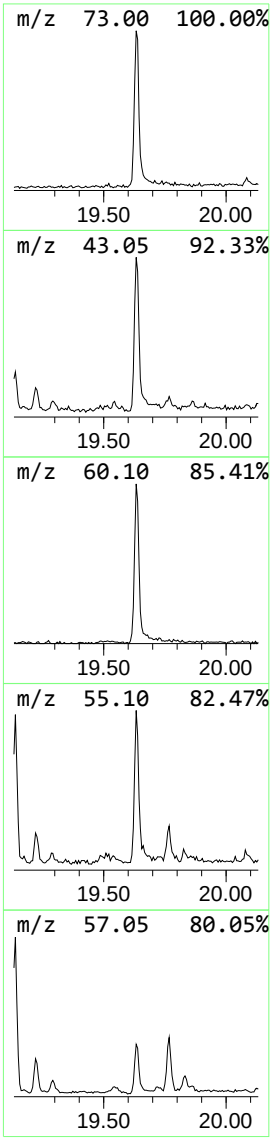
****

Peak Number 5 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.632	12.62 ng	623316	Chrysene-d12	21.759

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
4	n-Decanoic acid	172	C10H20O2	000334-48-5	45
5	Undecane, 5-methyl-	170	C12H26	001632-70-8	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012721\
Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
0051-045-COMP-E-ELUTRIATE

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

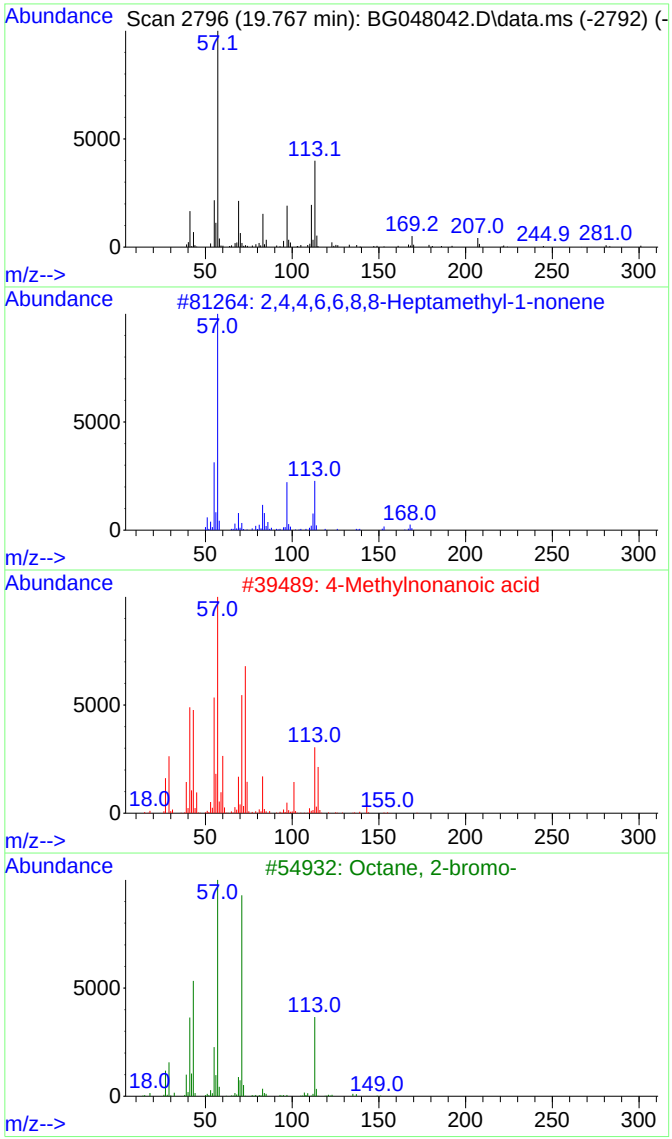
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

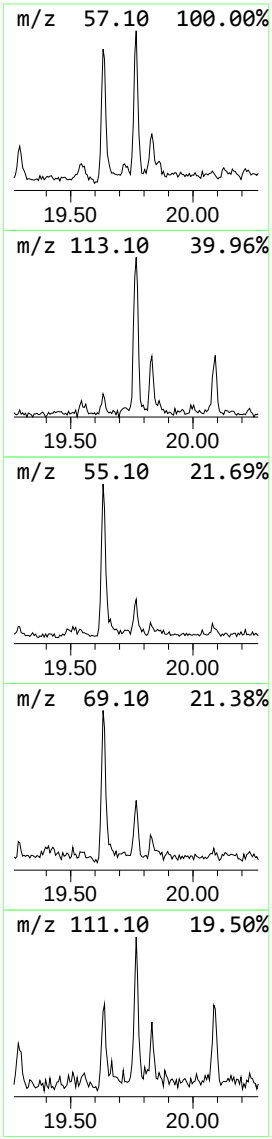
****

Peak Number 6 unknown19.767 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.767	2.62 ng	129408	Chrysene-d12	21.759

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	38
2	4-Methylnonanoic acid	172	C10H20O2	045019-28-1	25
3	Octane, 2-bromo-	192	C8H17Br	000557-35-7	23
4	2,4,4-Trimethyl-1-pentanol, hept...	326	C12H17F7O2	1000365-01-4	14
5	2,4,4-Trimethyl-1-pentanol, trif...	226	C10H17F3O2	1000365-19-5	14





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012721\
Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
0051-045-COMP-E-ELUTRIATE

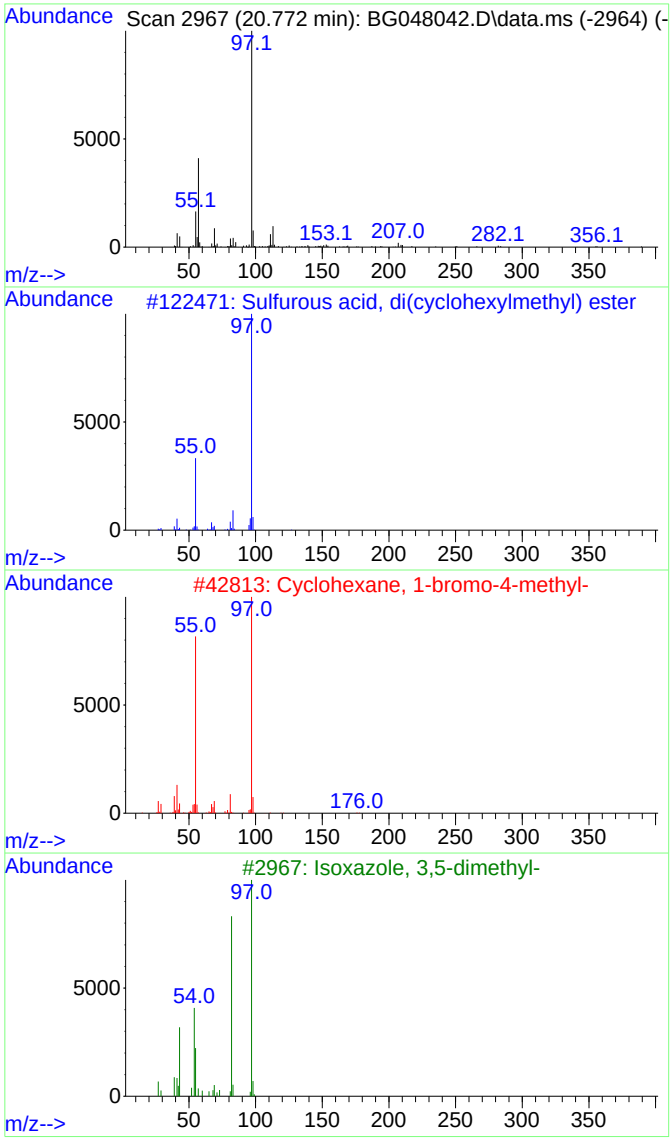
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

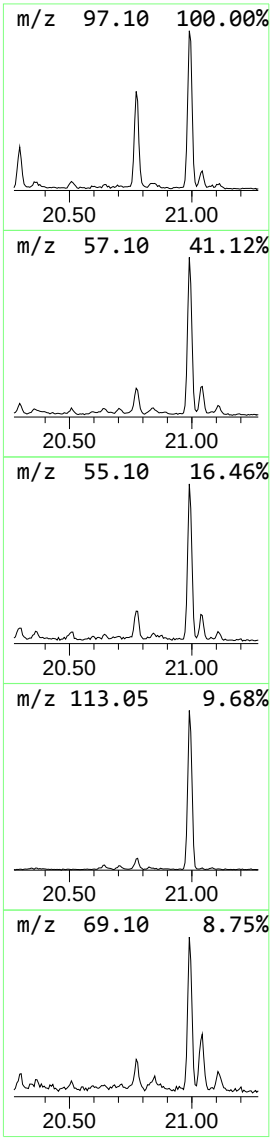
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Sulfurous acid, di(cyclohex... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.772	4.13 ng	203843	Chrysene-d12	21.759

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfurous acid, di(cyclohexylmet...	274	C14H26O3S	1000309-22-7	53
2	Cyclohexane, 1-bromo-4-methyl-	176	C7H13Br	006294-40-2	53
3	Isoxazole, 3,5-dimethyl-	97	C5H7NO	000300-87-8	42
4	2-Hexene, 3,4,4-trimethyl-	126	C9H18	053941-19-8	38
5	Cycloheptane, bromo-	176	C7H13Br	002404-35-5	36





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012721\
Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
0051-045-COMP-E-ELUTRIATE

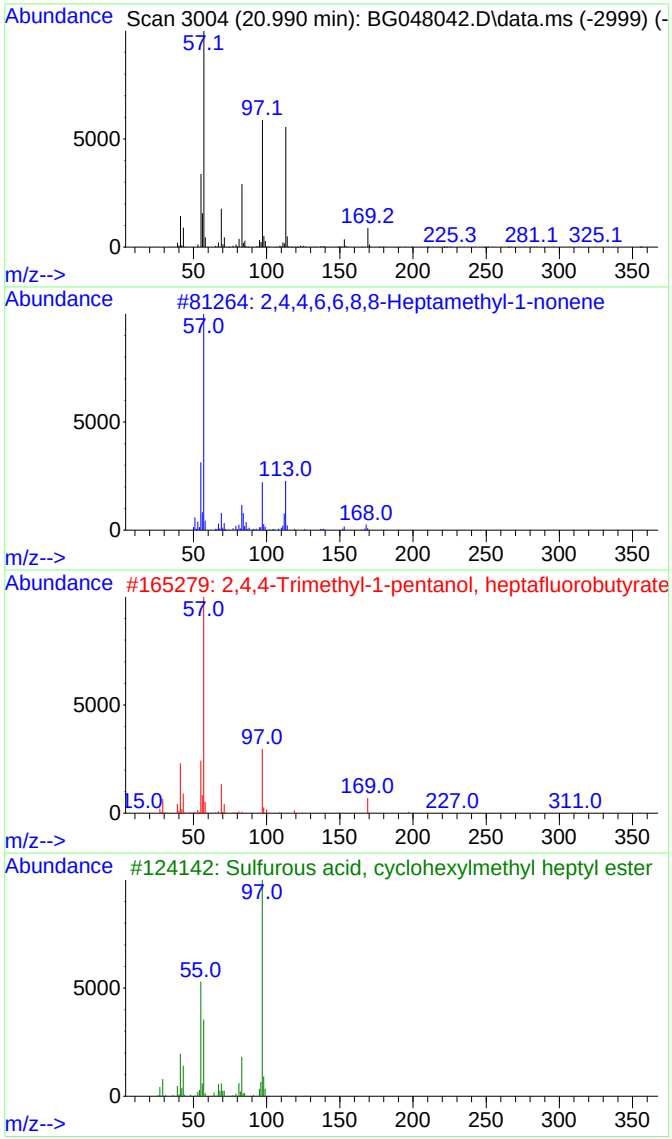
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

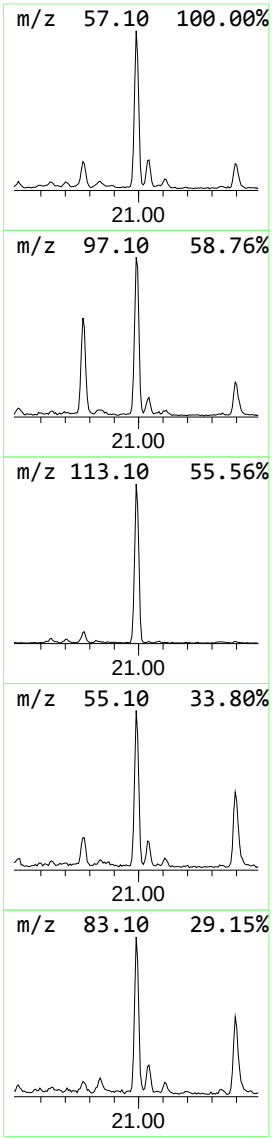
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 2,4,4,6,6,8,8-Heptamethyl-1... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.990	16.80 ng	829468	Chrysene-d12	21.759

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	50
2	2,4,4-Trimethyl-1-pentanol, hept...	326	C12H17F7O2	1000365-01-4	38
3	Sulfurous acid, cyclohexylmethyl...	276	C14H28O3S	1000309-21-7	37
4	Octane, 2-bromo-	192	C8H17Br	000557-35-7	35
5	Sulfurous acid, cyclohexylmethyl...	234	C11H22O3S	1000309-21-3	32





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012721\
Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
0051-045-COMP-E-ELUTRIATE

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

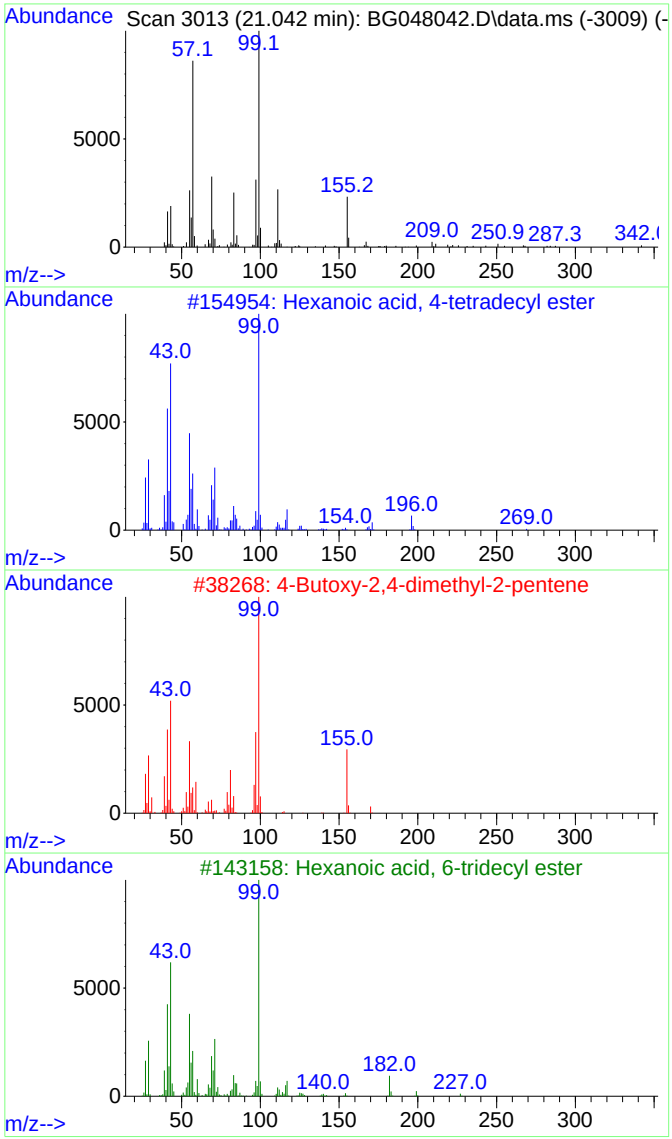
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

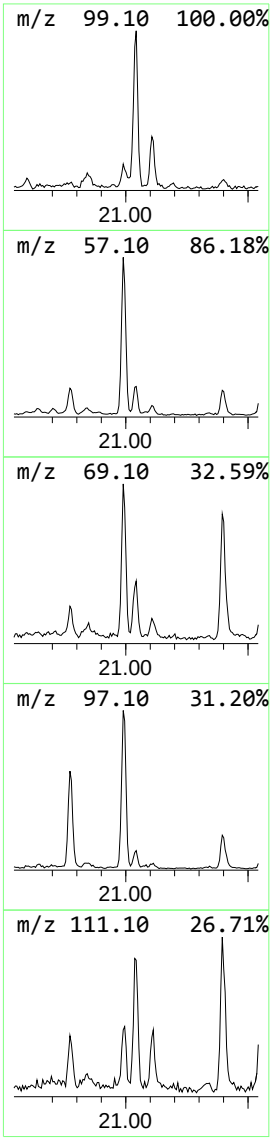
****

Peak Number 9 unknown21.043 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.043	4.24 ng	209275	Chrysene-d12	21.759

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 4-tetradecyl ester	312	C20H40O2	1000279-29-8	43
2	4-Butoxy-2,4-dimethyl-2-pentene	170	C11H22O	1000159-08-7	42
3	Hexanoic acid, 6-tridecyl ester	298	C19H38O2	1000279-29-5	38
4	Hexanoic acid, 3-chloroprop-2-en...	190	C9H15ClO2	1000299-36-1	38
5	Hexanoic acid, 4-tridecyl ester	298	C19H38O2	1000279-29-3	37





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012721\
Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
0051-045-COMP-E-ELUTRIATE

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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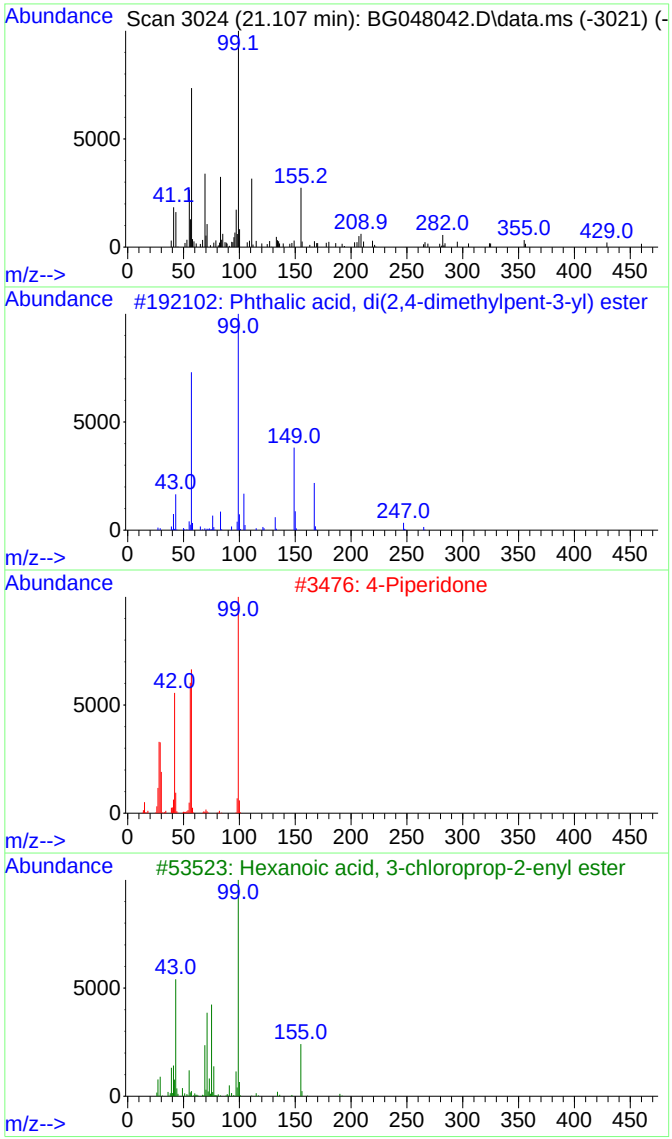
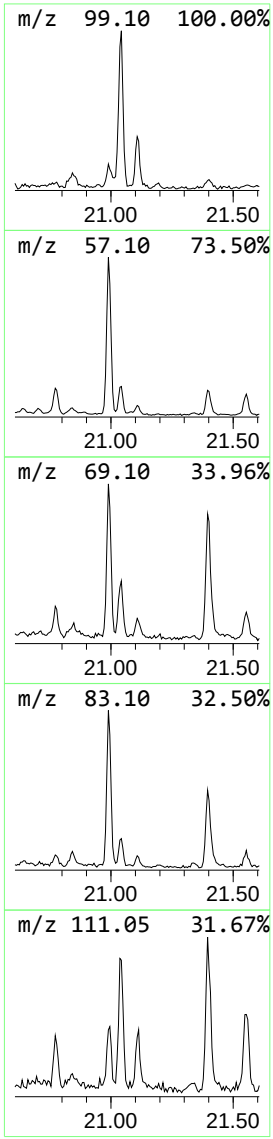
Peak Number 10 unknown21.107 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.107	2.01 ng	99303	Chrysene-d12	21.759

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phthalic acid, di(2,4-dimethylpe...	362	C22H34O4	1000356-85-5	43
2	4-Piperidone	99	C5H9NO	041661-47-6	43
3	Hexanoic acid, 3-chloroprop-2-en...	190	C9H15ClO2	1000299-36-1	38
4	1-Pyrrolidinocarbonylmethyl pipe...	197	C10H19N3O	039890-45-4	37
5	2H-Pyran-2-one, tetrahydro-6-nonyl-	226	C14H26O2	002721-22-4	37





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Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
0051-045-COMP-E-ELUTRIATE

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
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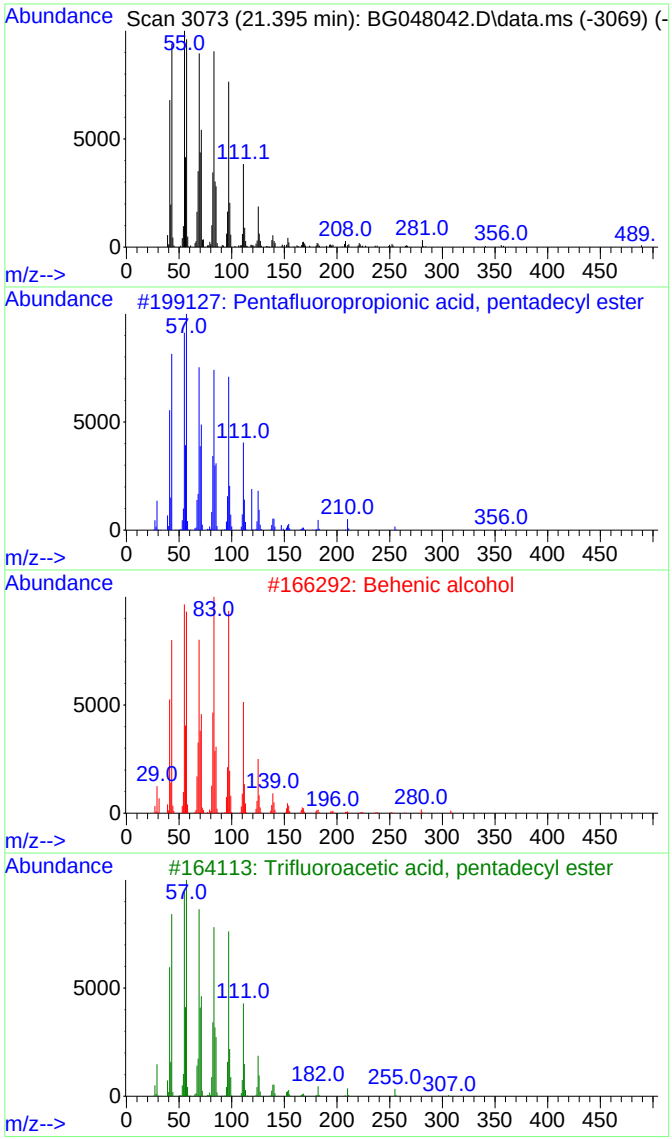
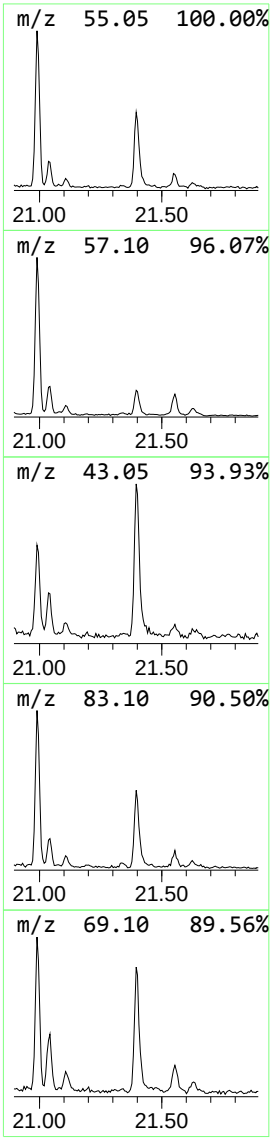
Peak Number 11 Pentafluoropropionic acid, ... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.395	9.27 ng	457835	Chrysene-d12	21.759

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
2	Behenic alcohol	326	C22H46O	000661-19-8	94
3	Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	94
4	1-Nonadecene	266	C19H38	018435-45-5	91
5	1-Heneicosyl formate	340	C22H44O2	077899-03-7	91





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Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : 0051-045-COMP-E-ELUTRIATE

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
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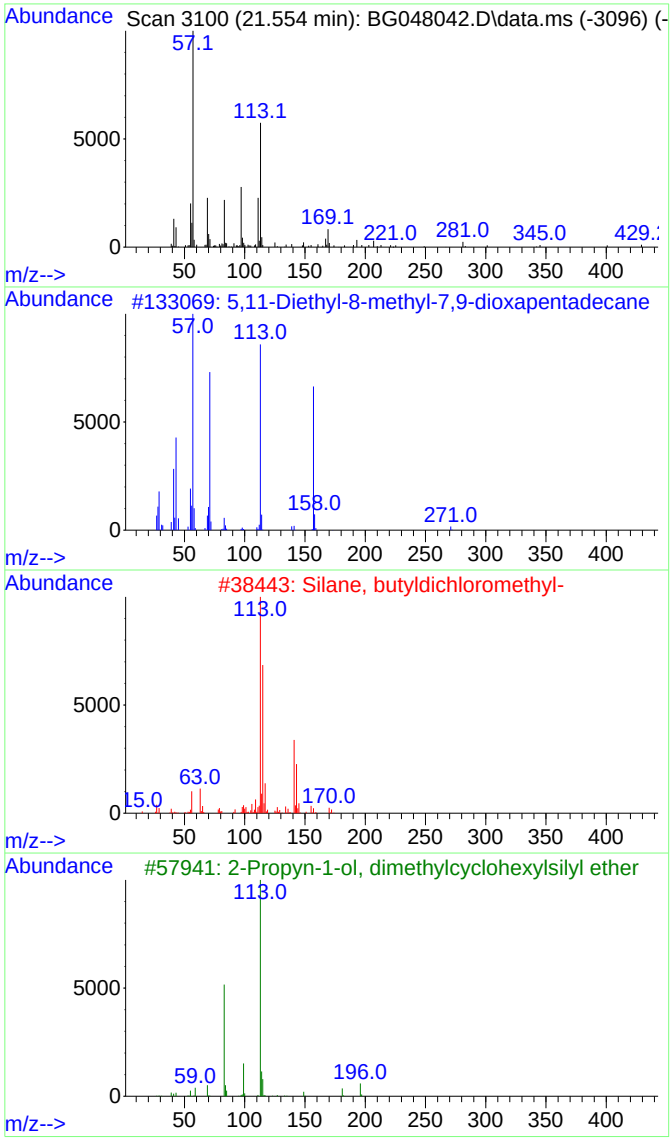
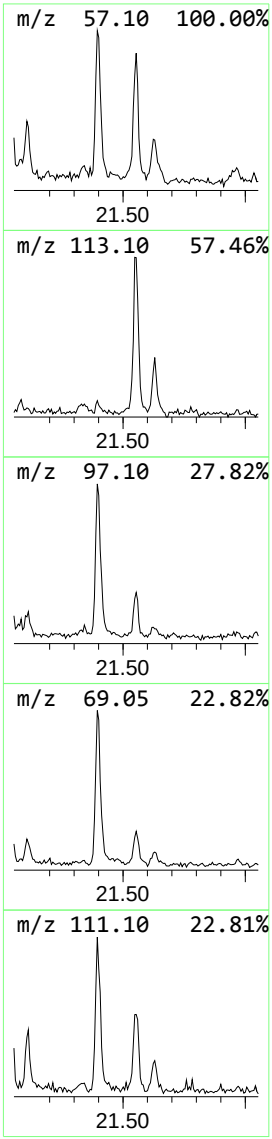
Peak Number 12 unknown21.554 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.554	2.06 ng	101949	Chrysene-d12	21.759

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,11-Diethyl-8-methyl-7,9-dioxap...	286	C18H38O2	1000334-83-1	32
2	Silane, butyldichloromethyl-	170	C5H12Cl2Si	018147-23-4	22
3	2-Propyn-1-ol, dimethylcyclohexy...	196	C11H20OSi	1000325-61-6	17
4	2,4,4-Trimethyl-1-pentanol, pent...	276	C11H17F5O2	1000365-51-0	14
5	2,2-Dimethyl-3-pentanol, heptafl...	312	C11H15F7O2	1000364-14-4	14





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Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
Sample : M1226-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
0051-045-COMP-E-ELUTRIATE

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
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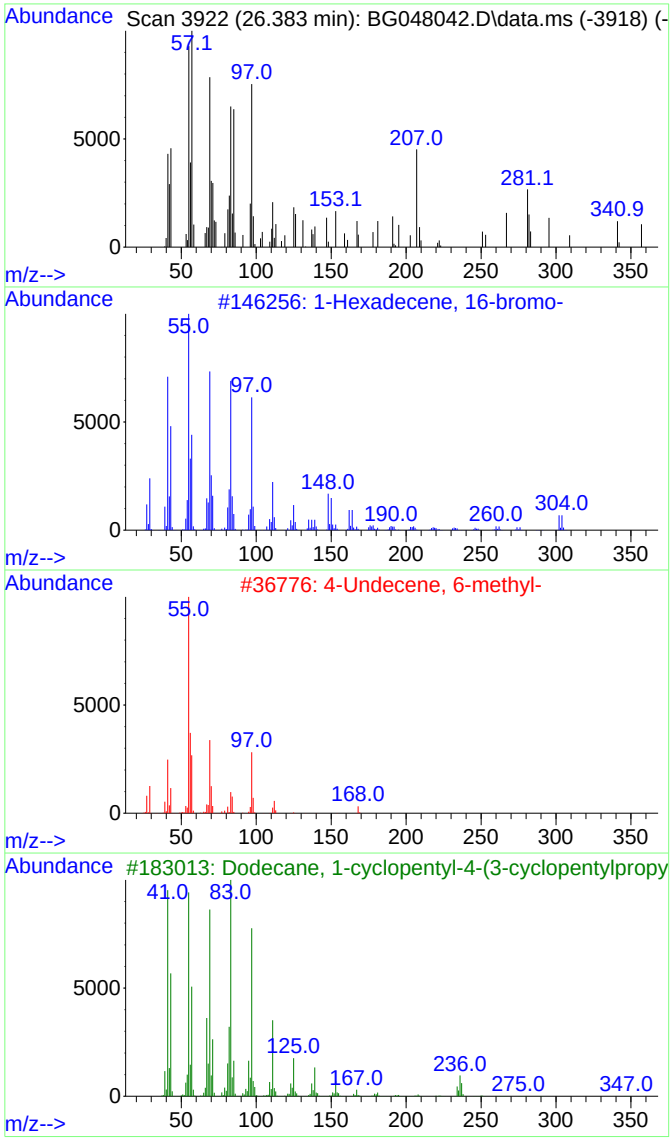
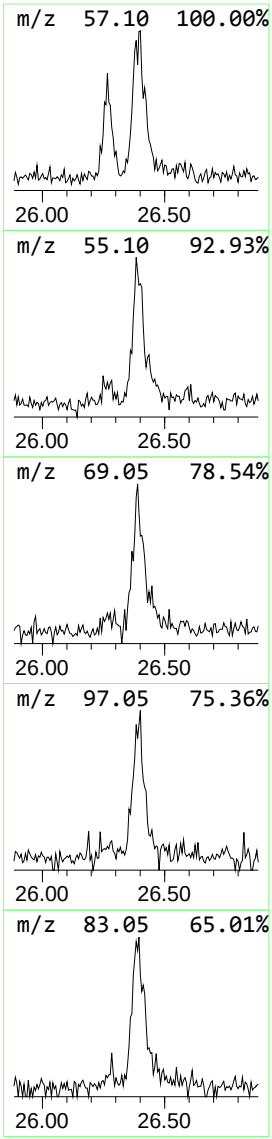
Peak Number 13 unknown26.383 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.383	2.54 ng	114824	Perylene-d12	25.049

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexadecene, 16-bromo-	302	C16H31Br	118625-56-2	43
2	4-Undecene, 6-methyl-	168	C12H24	1000061-85-4	43
3	Dodecane, 1-cyclopentyl-4-(3-cyc...	348	C25H48	007225-68-5	35
4	Octatriacontyl trifluoroacetate	647	C40H77F3O2	1000351-88-7	22
5	Carbonic acid, octadecyl 2,2,2-t...	444	C21H39Cl3O3	1000314-56-3	22





Tentatively Identified Compound (LSC) summary

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Data File : BG048042.D
Acq On : 28 Jan 2021 00:20
Operator : CG/JU
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ClientSampleId : 0051-045-COMP-E-ELUTRIATE

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
1,2-Benzenedica...	17.764	32.6	ng	1253910	4	17.488	769731	20.0
n-Hexadecanoic ...	18.369	20.0	ng	770146	4	17.488	769731	20.0
unknown19.139	19.139	9.7	ng	371672	4	17.488	769731	20.0
unknown19.221	19.221	3.1	ng	120853	4	17.488	769731	20.0
Octadecanoic acid	19.632	12.6	ng	623316	5	21.759	987511	20.0
unknown19.767	19.767	2.6	ng	129408	5	21.759	987511	20.0
Sulfurous acid,...	20.772	4.1	ng	203843	5	21.759	987511	20.0
2,4,4,6,6,8,8-H...	20.990	16.8	ng	829468	5	21.759	987511	20.0
unknown21.043	21.043	4.2	ng	209275	5	21.759	987511	20.0
unknown21.107	21.107	2.0	ng	99303	5	21.759	987511	20.0
Pentafluoroprop...	21.395	9.3	ng	457835	5	21.759	987511	20.0
unknown21.554	21.554	2.1	ng	101949	5	21.759	987511	20.0
unknown26.383	26.383	2.5	ng	114824	6	25.049	904730	20.0