LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060421\
Data File : BF124138.D
Acq On : 4 Jun 2021 14:16
Operator : JU/CG
Sample : M2587-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : FK-GF-02-073-A

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_F\METHODS\8270-BF060321.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF124138.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.152	3	11	17	rVB	500439	637749	53.37%	5.886%
2	2.257	25	29	35	rBB	28675	36530	3.06%	0.337%
3	5.093	507	511	516	rVB	33235	38994	3.26%	0.360%
4	5.493	575	579	583	rBV	1110243	1089393	91.17%	10.054%
5	6.504	746	751	754	rBV	1081484	1054596	88.26%	9.733%
6	6.863	809	812	816	rVB	378627	298480	24.98%	2.755%
7	7.428	903	908	911	rBV	852964	722920	60.50%	6.672%
8	8.045	1010	1013	1018	rBV	72321	66663	5.58%	0.615%
9	8.145	1023	1030	1033	rBV	441355	366275	30.65%	3.380%
10	9.222	1209	1213	1216	rBV	1509575	1194891	100.00%	11.028%
11	9.298	1223	1226	1231	rBV4	22881	29814	2.50%	0.275%
12	9.616	1277	1280	1283	rBV3	13424	17133	1.43%	0.158%
13	9.904	1325	1329	1332	rBV	436599	365805	30.61%	3.376%
14	9.933	1332	1334	1337	rVB2	22899	16989	1.42%	0.157%
15	10.328	1399	1401	1405	rVB	17791	15830	1.32%	0.146%
16	10.445	1417	1421	1425	rVB3	22245	26541	2.22%	0.245%
17	10.692	1459	1463	1466	rBV	726987	632931	52.97%	5.842%
18	10.816	1479	1484	1487	rBV3	22003	30430	2.55%	0.281%
19	11.251	1554	1558	1560	rBV4	20904	19425	1.63%	0.179%
20	11.286	1560	1564	1568	rVB3	20227	20781	1.74%	0.192%
21	11.386	1578	1581	1584	rBV	361891	330440	27.65%	3.050%
22	11.410	1584	1585	1589	rVV	150726	100287	8.39%	0.926%
23	11.463	1592	1594	1597	rVB	36198	31009	2.60%	0.286%
24	11.616	1618	1620	1625	rVB3	15088	16236	1.36%	0.150%
25	11.680	1628	1631	1634	rBV3	19006	18446	1.54%	0.170%
26	11.780	1644	1648	1650	rBV	83146	66691	5.58%	0.616%
27	11.916	1668	1671	1674	rVB2	61060	57254	4.79%	0.528%
28	12.004	1682	1686	1688	rBV	32634	31677	2.65%	0.292%
29	12.186	1714	1717	1719	rBV3	20633	17202	1.44%	0.159%
30	12.210	1719	1721	1724	rVB4	19552	17978	1.50%	0.166%
31	12.445	1757	1761	1762	rBV4	20008	28445	2.38%	0.263%
32	12.469	1762	1765	1767	rVV	309658	320347	26.81%	2.957%
33	12.486	1767	1768	1773	rVB2	229220	162371	13.59%	1.499%
34	12.574	1779	1783	1785	rBV2	41902	41440	3.47%	0.382%
35	12.604	1785	1788	1791	rVV	185020	162547	13.60%	1.500%
36	12.698	1801	1804	1807	rBV5	25046	27034	2.26%	0.250%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060421\
Data File : BF124138.D
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Operator : JU/CG
Sample : M2587-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : FK-GF-02-073-A

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_F\METHODS\8270-BF060321.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	12.833	1822	1827	1832	rVV	163573	173774	14.54%	1.604%
38	12.874	1832	1834	1840	rVB6	17676	24241	2.03%	0.224%
39	12.974	1847	1851	1858	rVB	1163181	994166	83.20%	9.176%
40	13.139	1876	1879	1880	rBV3	18144	19679	1.65%	0.182%
41	13.163	1880	1883	1886	rVB2	35968	36732	3.07%	0.339%
42	13.204	1886	1890	1891	rBV4	28824	24294	2.03%	0.224%
43	13.227	1891	1894	1896	rVB4	28048	29404	2.46%	0.271%
44	13.263	1896	1900	1902	rBV5	26482	37392	3.13%	0.345%
45	13.668	1966	1969	1971	rBV4	25861	21736	1.82%	0.201%
46	13.880	2001	2005	2011	rVB6	66666	94980	7.95%	0.877%
47	14.033	2026	2031	2033	rBV2	343473	392642	32.86%	3.624%
48	14.057	2033	2035	2041	rVB	87806	77436	6.48%	0.715%
49	14.133	2046	2048	2051	rBV4	18347	14545	1.22%	0.134%
50	14.498	2108	2110	2112	rBV2	36086	39104	3.27%	0.361%
51	15.080	2206	2209	2216	rVB2	74246	134703	11.27%	1.243%
52	15.380	2258	2260	2267	rVB3	50032	71282	5.97%	0.658%
53	15.451	2269	2272	2278	rVB	58151	82127	6.87%	0.758%
54	15.515	2278	2283	2286	rBV	200673	251278	21.03%	2.319%
55	16.227	2402	2404	2408	rVB5	42256	54736	4.58%	0.505%
56	16.662	2476	2478	2485	rVB8	32978	60327	5.05%	0.557%
57	17.092	2548	2551	2552	rBV3	27997	28789	2.41%	0.266%
58	17.339	2591	2593	2596	rVB4	23894	17743	1.48%	0.164%
59	17.439	2608	2610	2619	rVB10	21224	42289	3.54%	0.390%

Sum of corrected areas: 10834973

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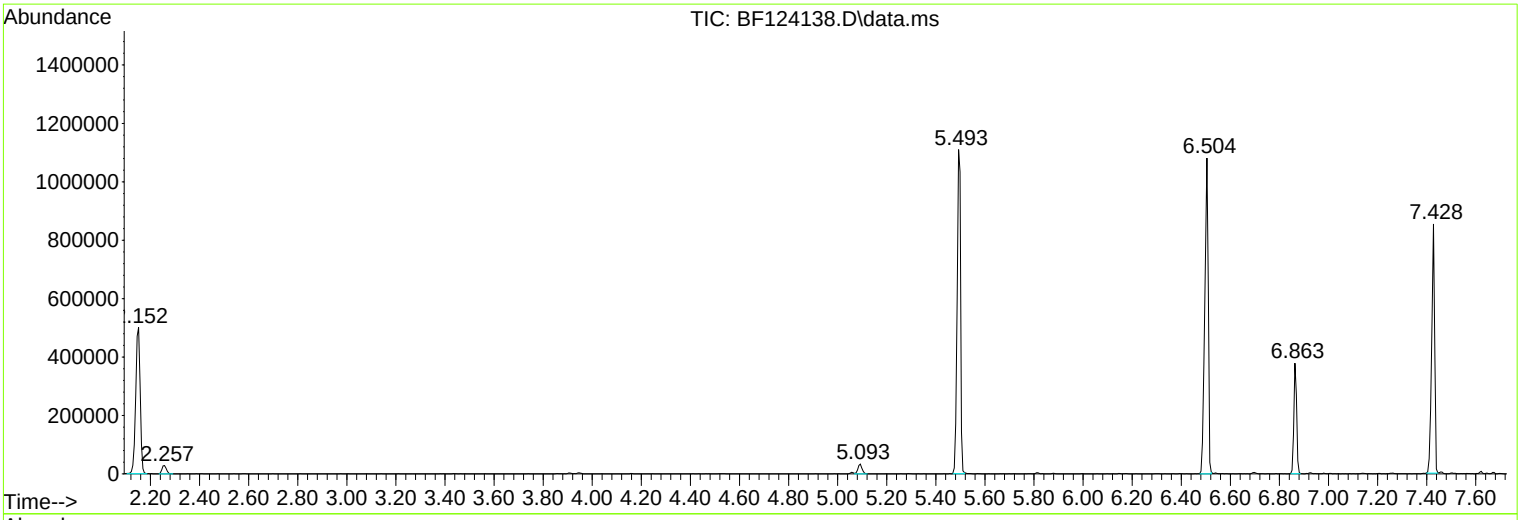
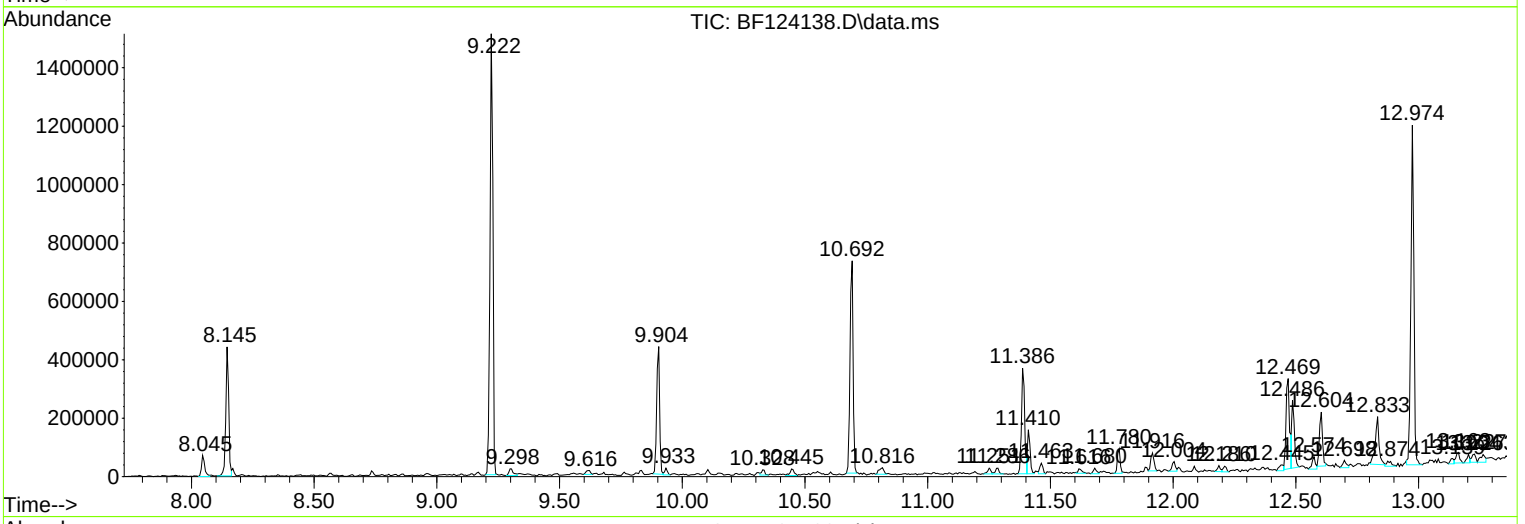
Misc :
ALS Vial : 11 Sample Multiplier: 1

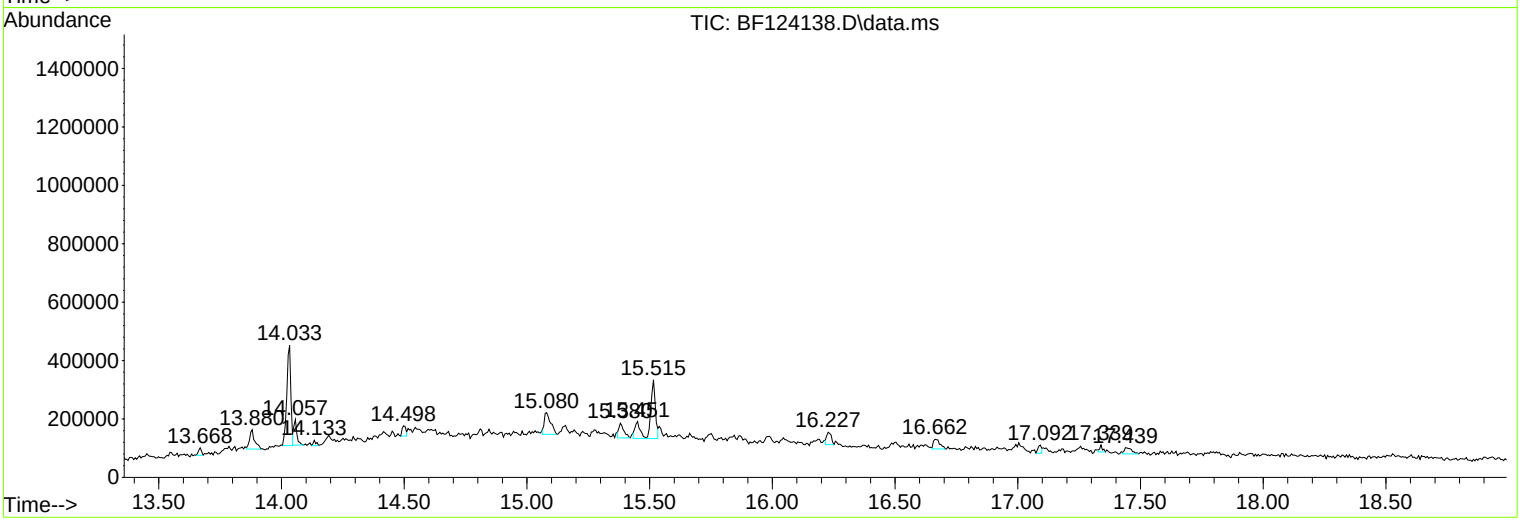
Instrument :
BNA_F
ClientSampleId :
FK-GF-02-073-A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.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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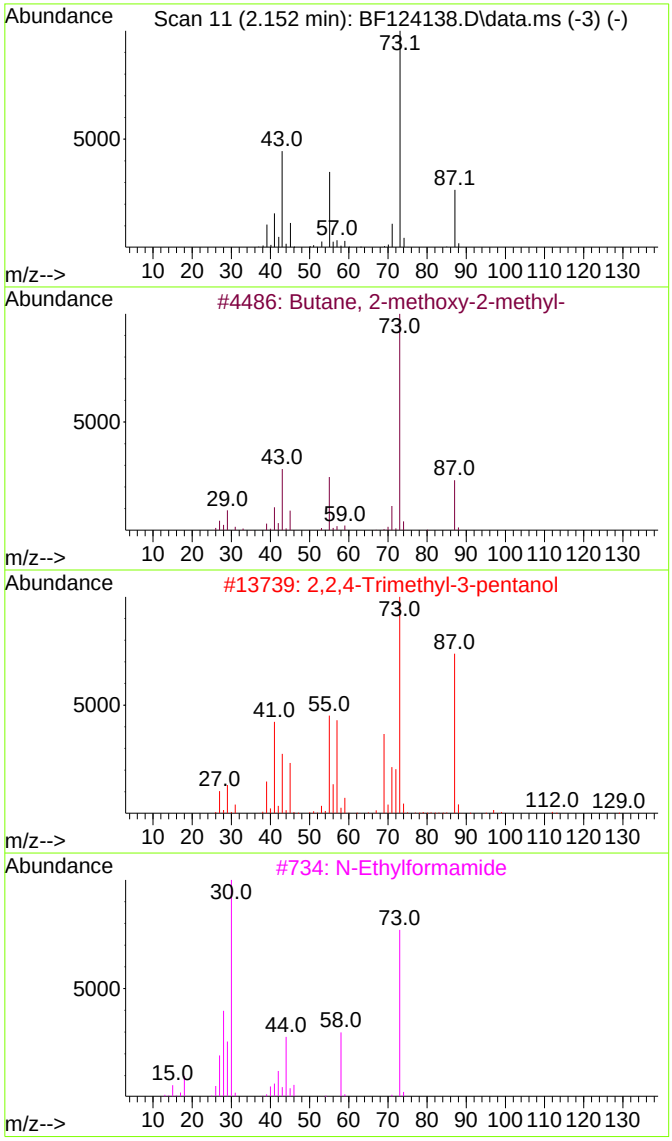
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

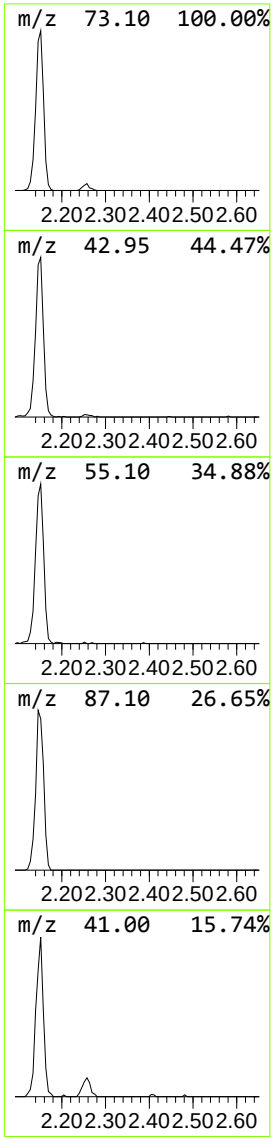
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.152	42.73 ng	637749	1,4-Dichlorobenzene-d4	6.863

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	38
3	N-Ethylformamide	73	C3H7NO	000627-45-2	9
4	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5	3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060421\
Data File : BF124138.D
Acq On : 4 Jun 2021 14:16
Operator : JU/CG
Sample : M2587-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : FK-GF-02-073-A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

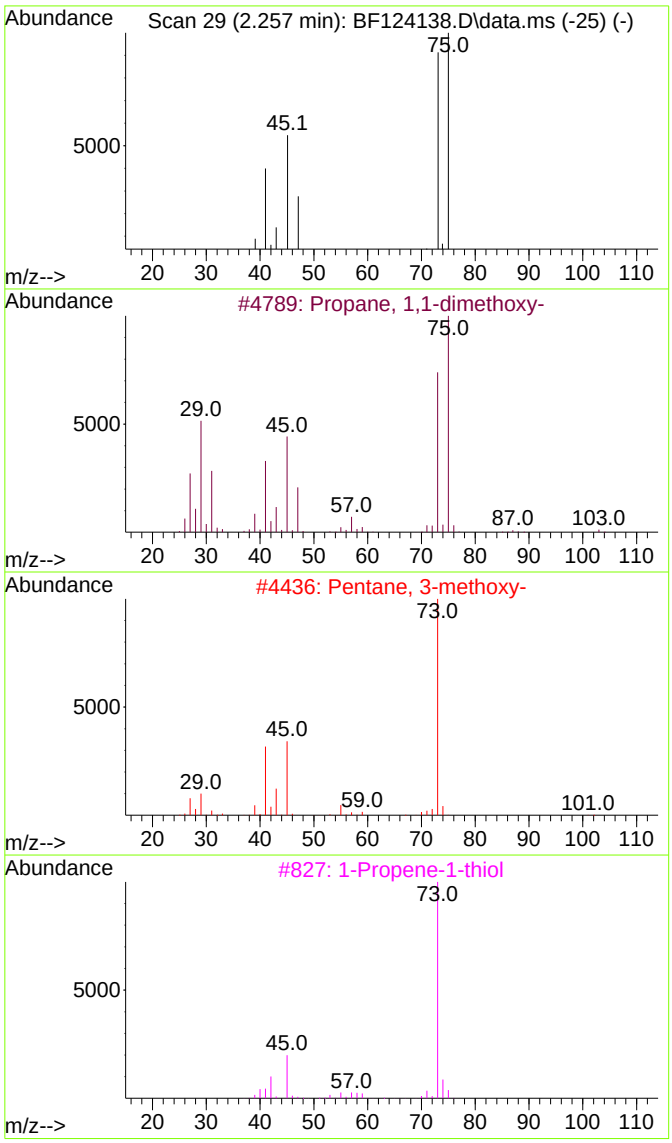
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

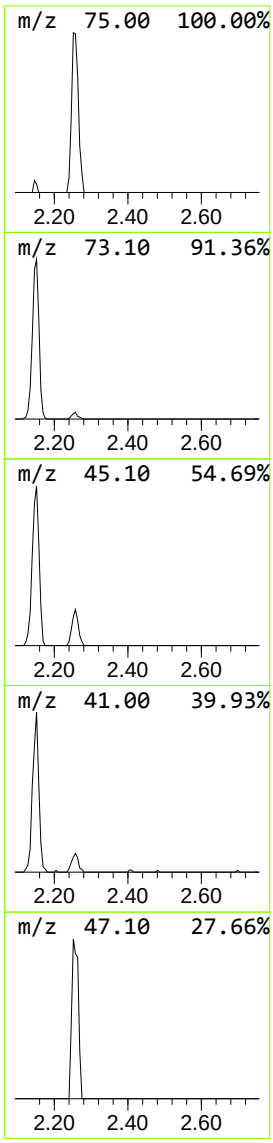
****

Peak Number 2 Propane, 1,1-dimethoxy- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.257	2.45 ng	36530	1,4-Dichlorobenzene-d4	6.863

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	74
2	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9
3	1-Propene-1-thiol	74	C3H6S	000925-89-3	7
4	Methane, nitroso-	45	CH3NO	000865-40-7	4
5	Methyl propyl ether	74	C4H10O	000557-17-5	4





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060421\
Data File : BF124138.D
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Operator : JU/CG
Sample : M2587-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FK-GF-02-073-A

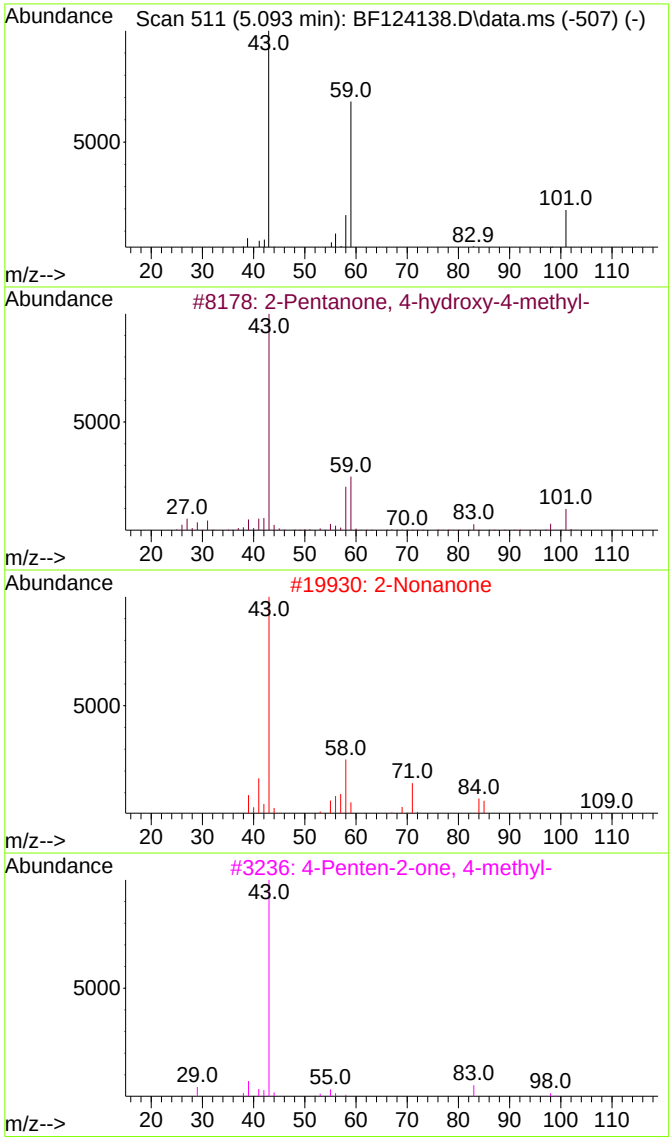
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

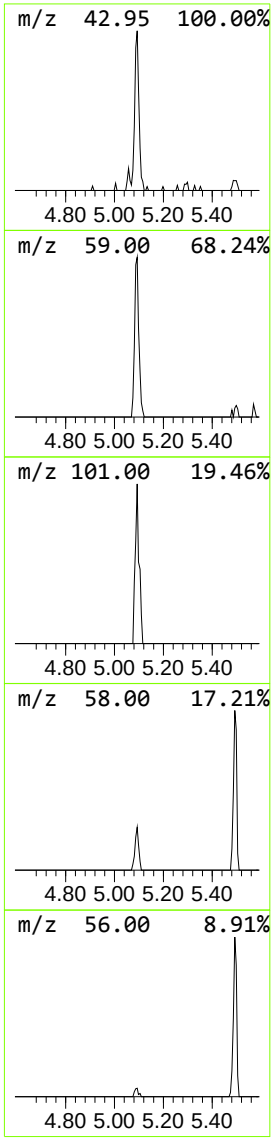
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.093	2.61 ng	38994	1,4-Dichlorobenzene-d4	6.863

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	2-Nonanone	142	C9H18O	000821-55-6	17
3	4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	10
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9
5	Acetone	58	C3H6O	000067-64-1	9





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Instrument :
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ClientSampleId :
FK-GF-02-073-A

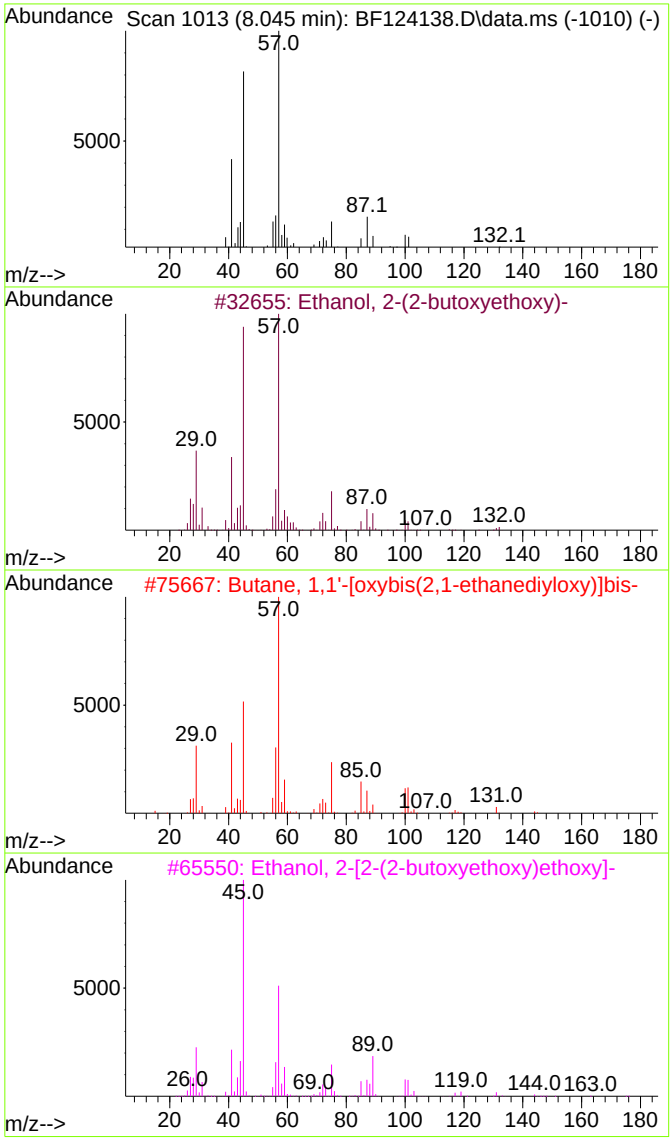
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

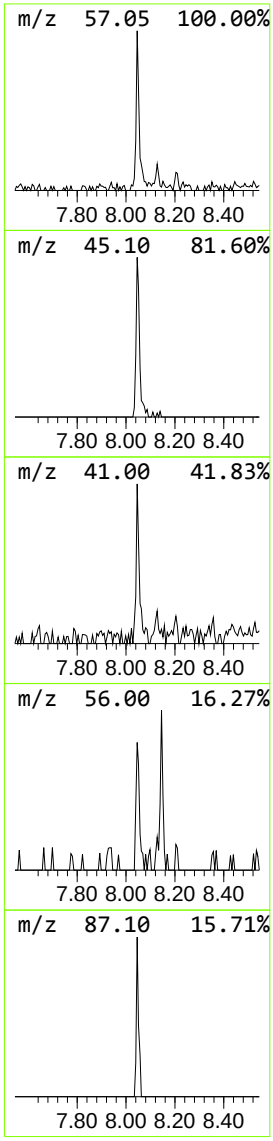
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.045	3.64 ng	66663	Naphthalene-d8	8.145

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Butane, 1,1'-[oxybis(2,1-ethaned...	218	C12H26O3	000112-73-2	83
3	Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	72
4	Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	53
5	Pentaethylene glycol	238	C10H22O6	004792-15-8	42





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060421\
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Operator : JU/CG
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ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FK-GF-02-073-A

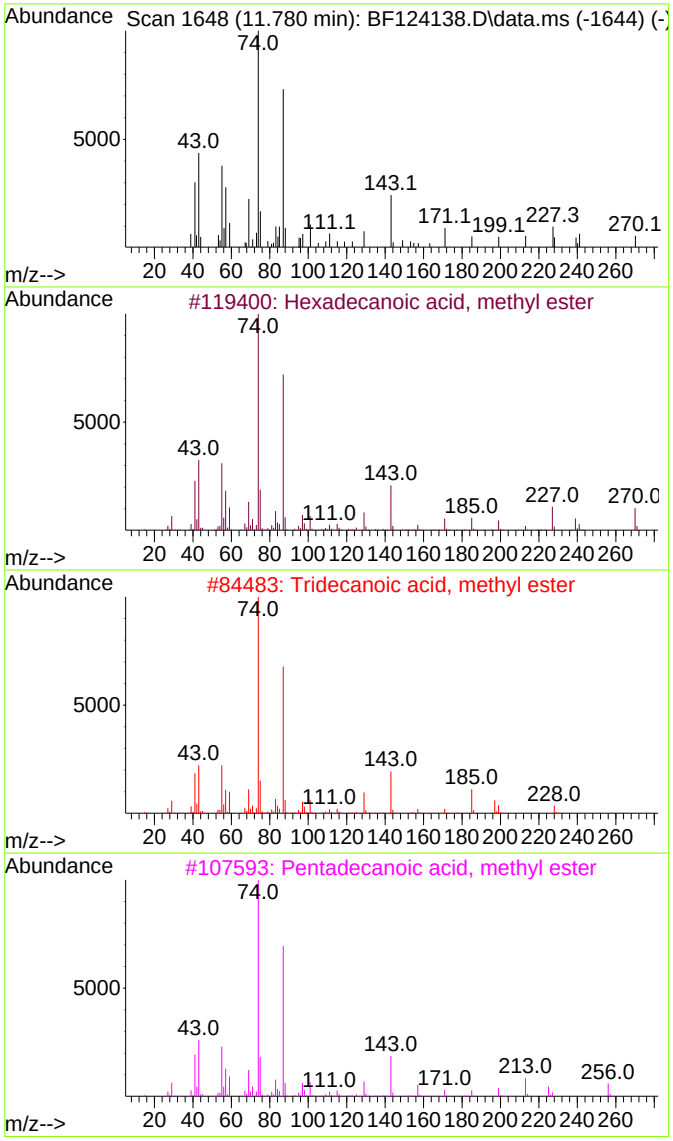
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

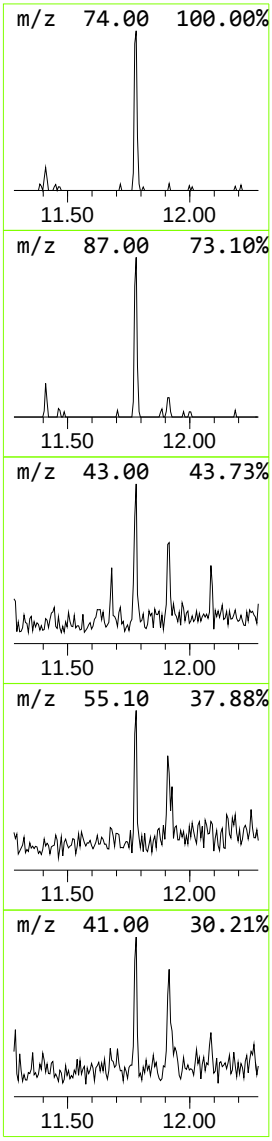
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Hexadecanoic acid, methyl e... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.780	4.04 ng	66691	Phenanthrene-d10	11.386

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	97
2	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	91
3	Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	72
4	Undecanoic acid, 10-methyl-, met...	214	C13H26O2	005129-56-6	72
5	Methyl 11-methyl-dodecanoate	228	C14H28O2	1000336-45-1	70





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Instrument : BNA_F
ClientSampleId : FK-GF-02-073-A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

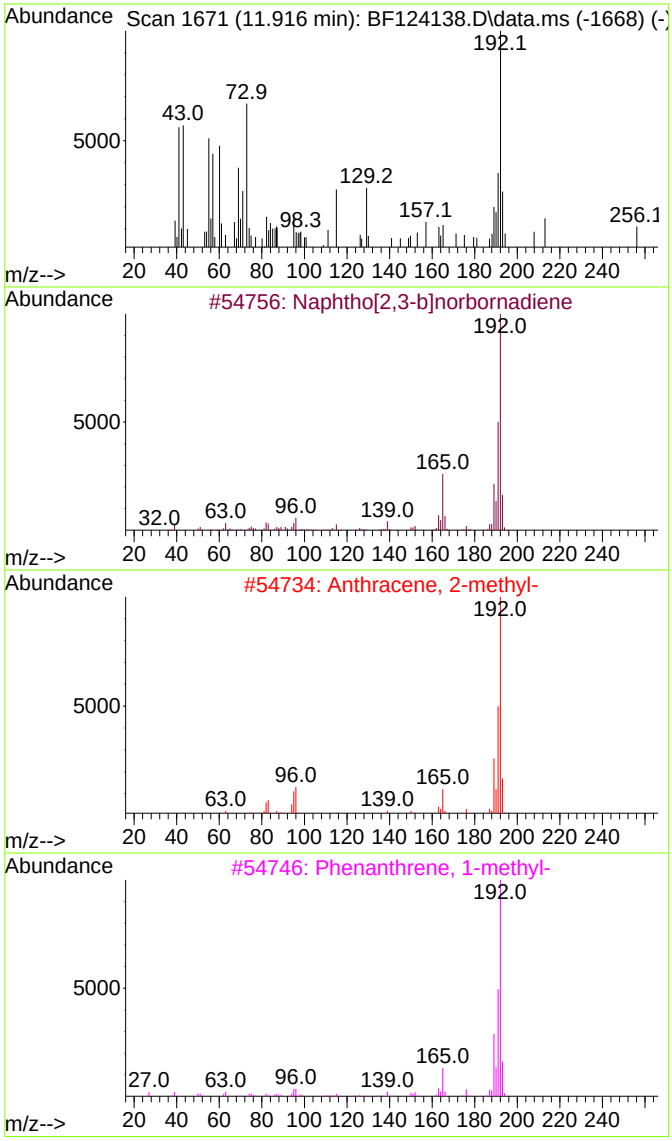
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

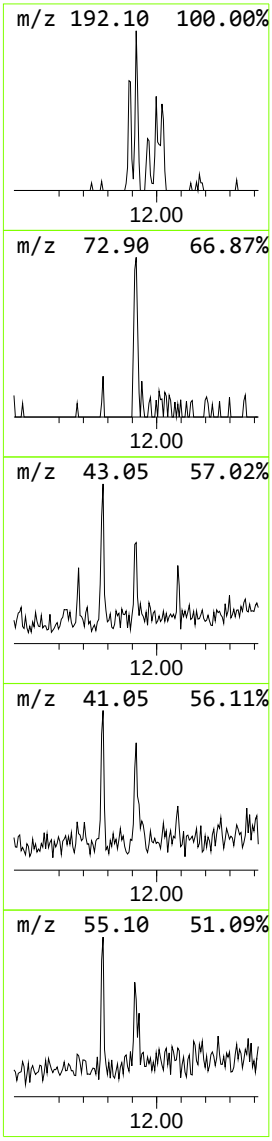
****

Peak Number 6 unknown11.916 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.916	3.47 ng	57254	Phenanthrene-d10	11.386

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	38
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	38
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	38
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	38
5	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060421\
Data File : BF124138.D
Acq On : 4 Jun 2021 14:16
Operator : JU/CG
Sample : M2587-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FK-GF-02-073-A

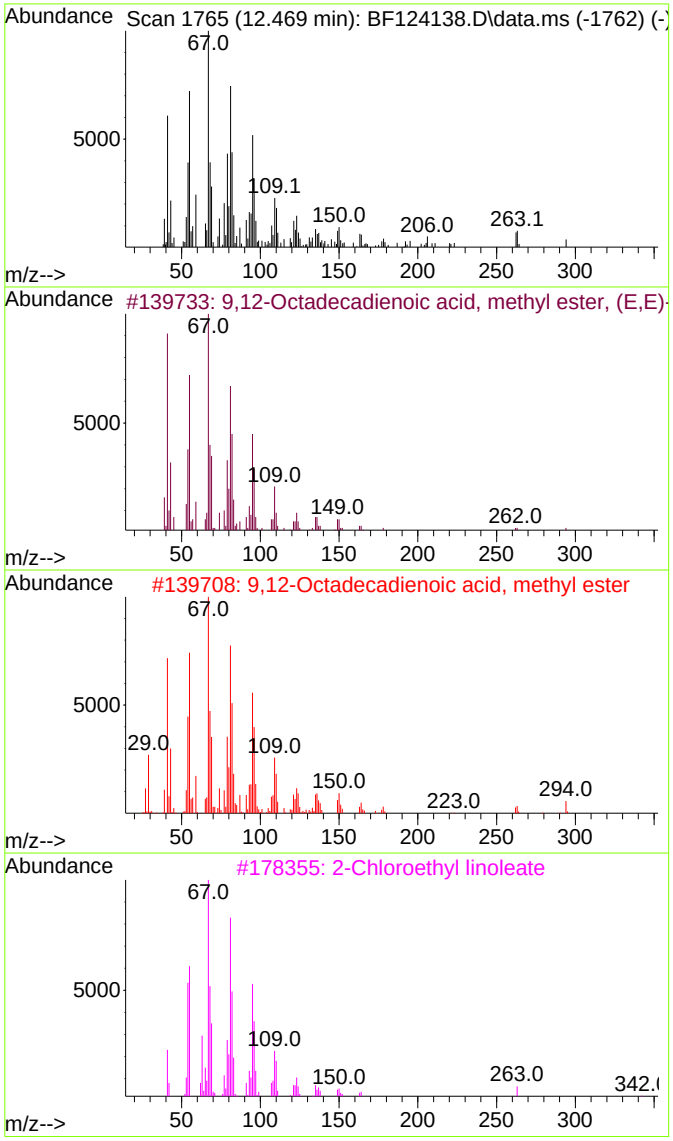
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

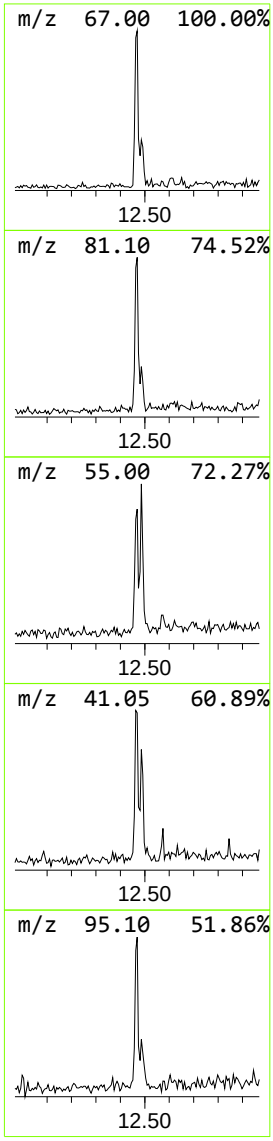
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 9,12-Octadecadienoic acid, ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.469	19.39 ng	320347	Phenanthrene-d10	11.386

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	98
2	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	98
3	2-Chloroethyl linoleate	342	C20H35ClO2	025525-76-2	93
4	Methyl 7,12-octadecadienoate	294	C19H34O2	1000336-43-8	90
5	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	89





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Data File : BF124138.D
Acq On : 4 Jun 2021 14:16
Operator : JU/CG
Sample : M2587-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FK-GF-02-073-A

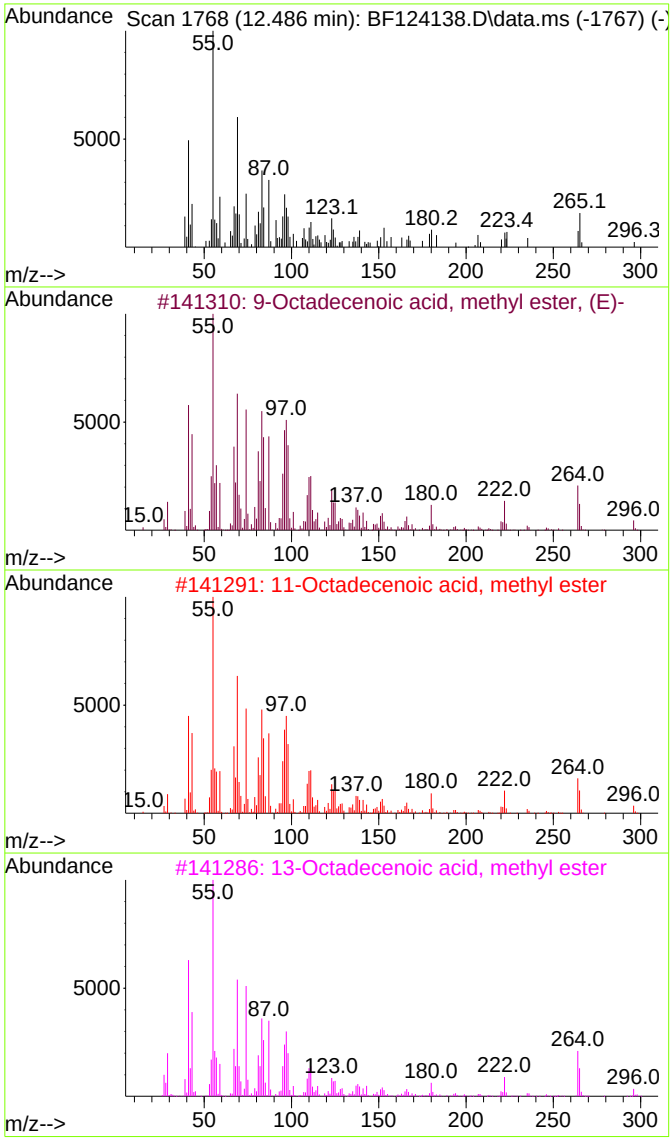
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

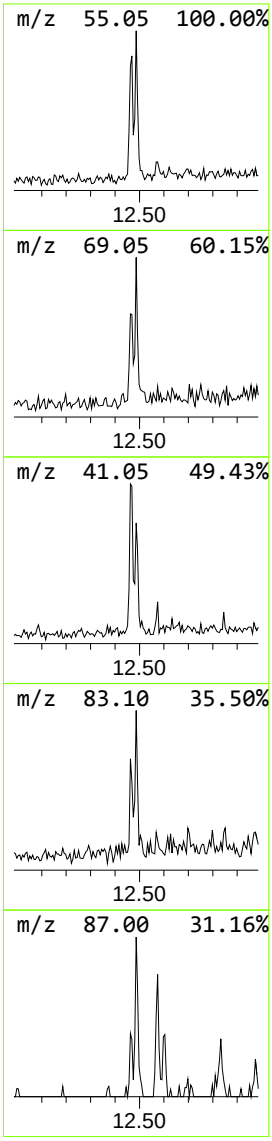
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 9-Octadecenoic acid, methyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.486	9.83 ng	162371	Phenanthrene-d10	11.386

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	95
2	11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	78
3	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	64
4	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	58
5	8-Octadecenoic acid, methyl ester	296	C19H36O2	002345-29-1	55





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Acq On : 4 Jun 2021 14:16
Operator : JU/CG
Sample : M2587-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : FK-GF-02-073-A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.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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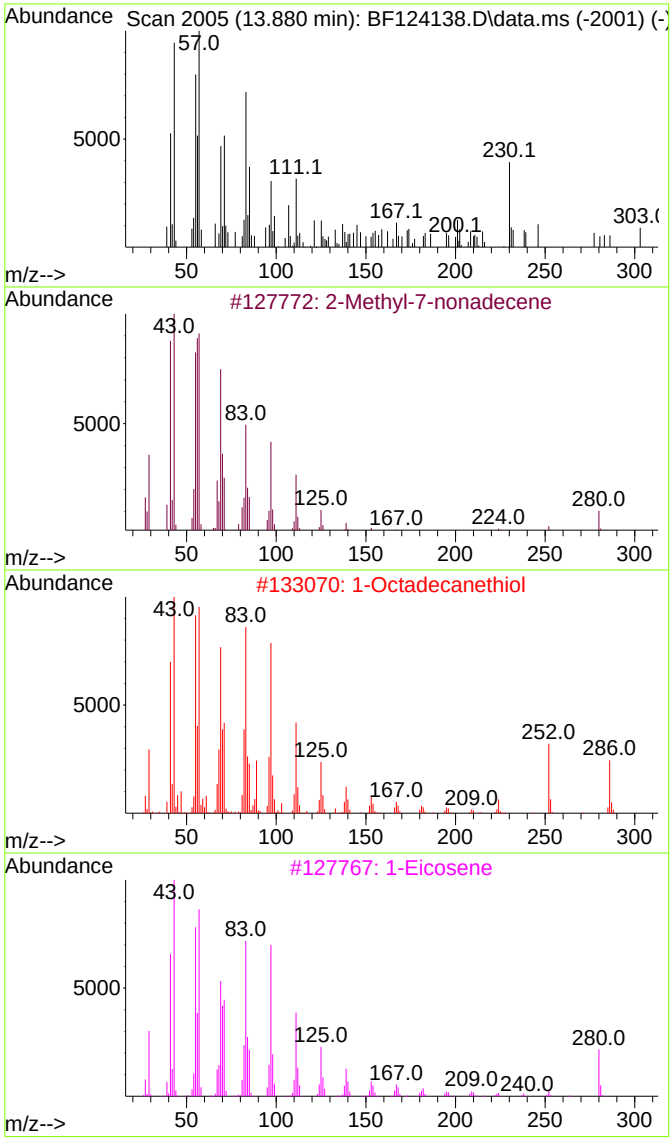
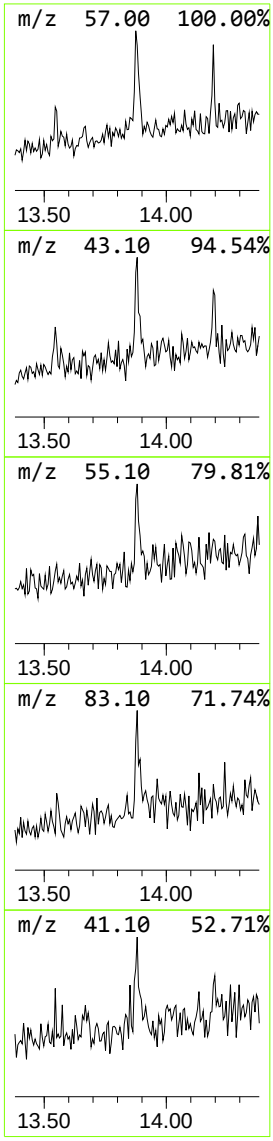
Peak Number 9 unknown13.880 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.880	4.84 ng	94980	Chrysene-d12	14.033

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-7-nonadecene	280	C20H40	219750-68-2	38
2	1-Octadecanethiol	286	C18H38S	002885-00-9	37
3	1-Eicosene	280	C20H40	003452-07-1	32
4	Cyclohexadecane, 1,2-diethyl-	280	C20H40	1000155-85-3	30
5	Tetradecanoic acid, 2-hydroxy-, ...	258	C15H30O3	056009-40-6	22





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Acq On : 4 Jun 2021 14:16
Operator : JU/CG
Sample : M2587-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : FK-GF-02-073-A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.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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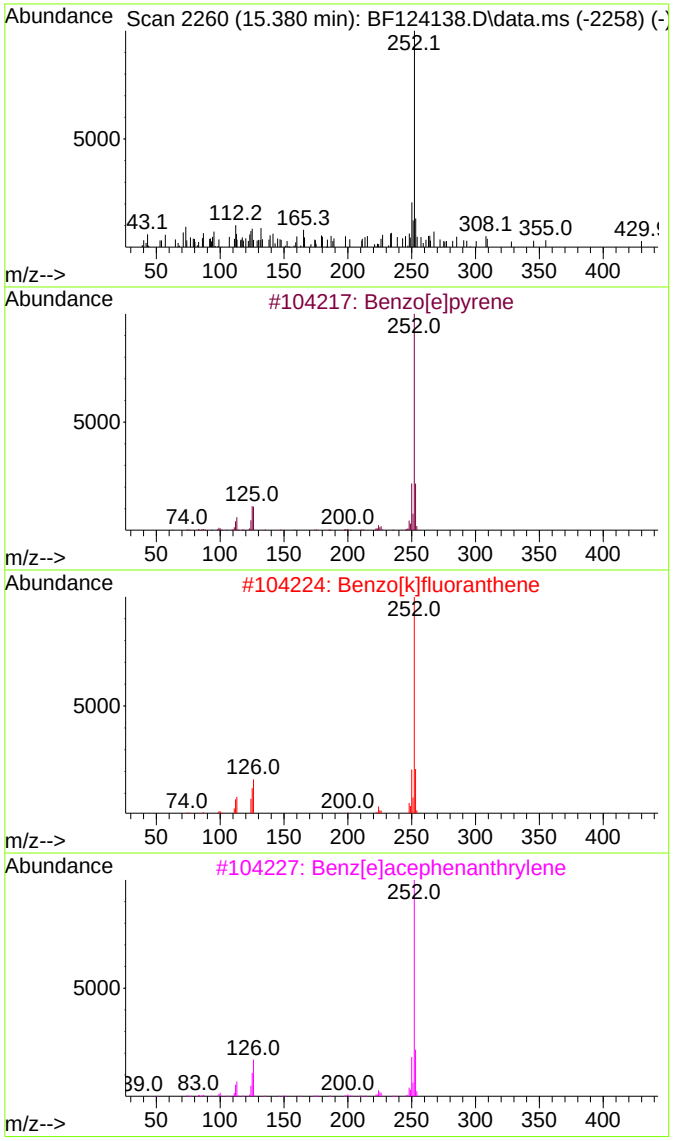
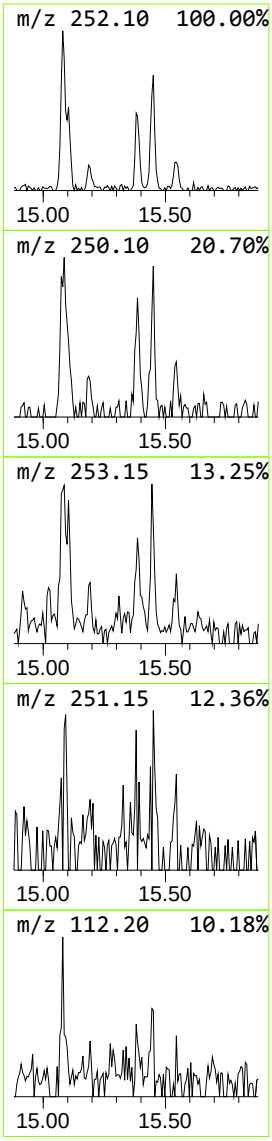
Peak Number 10 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.380	5.67 ng	71282	Perylene-d12	15.515

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	58
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	58
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	58
4	Benzo[a]pyrene	252	C20H12	000050-32-8	52
5	Perylene	252	C20H12	000198-55-0	52





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Acq On : 4 Jun 2021 14:16
Operator : JU/CG
Sample : M2587-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : FK-GF-02-073-A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.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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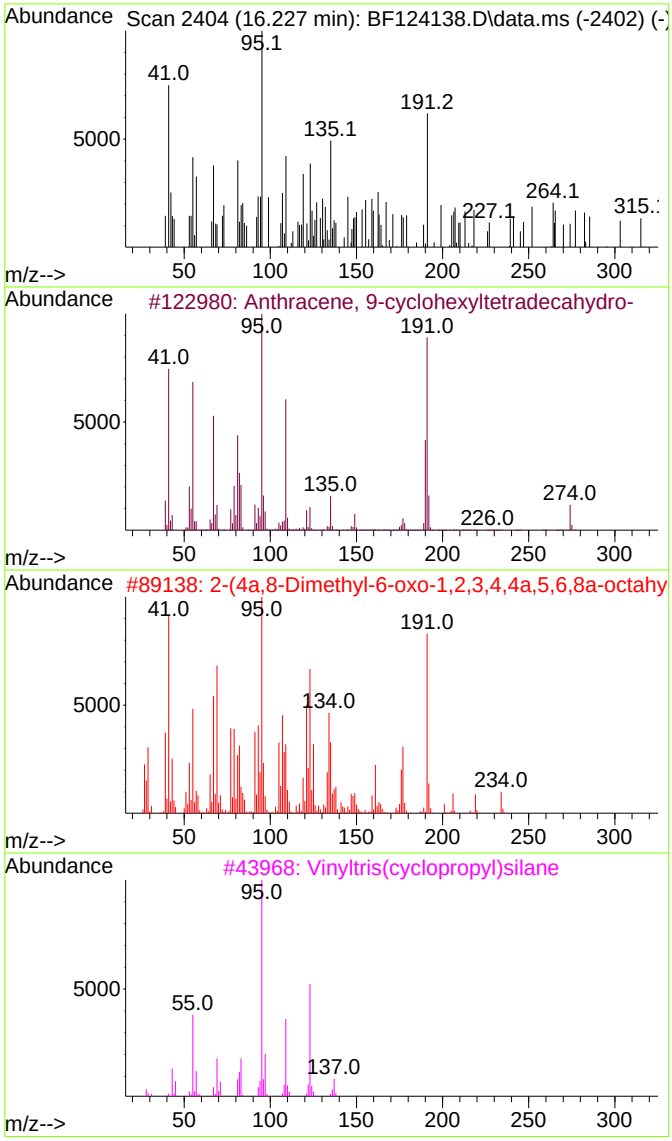
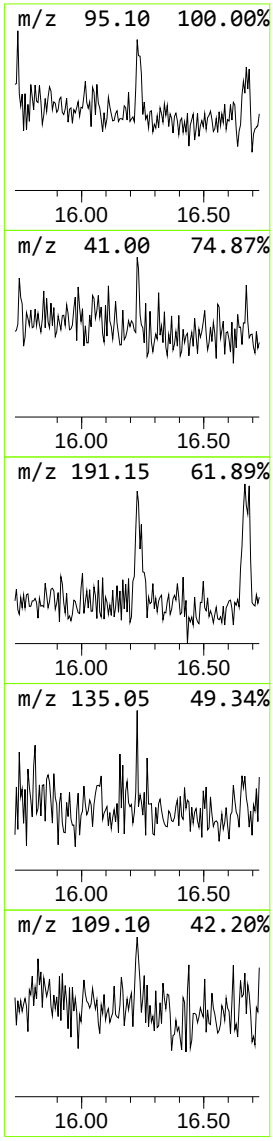
Peak Number 11 Anthracene, 9-cyclohexyltet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.227	4.36 ng	54736	Perylene-d12	15.515

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 9-cyclohexyltetradec...	274	C20H34	055255-70-4	55
2	2-(4a,8-Dimethyl-6-oxo-1,2,3,4,4...	234	C15H22O2	1000190-21-8	47
3	Vinyltris(cyclopropyl)silane	178	C11H18Si	1000109-97-3	35
4	Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	25
5	2,2,6-Trimethyl-1-(2-methyl-cycl...	218	C15H22O	1000188-72-8	25





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Acq On : 4 Jun 2021 14:16
Operator : JU/CG
Sample : M2587-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FK-GF-02-073-A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.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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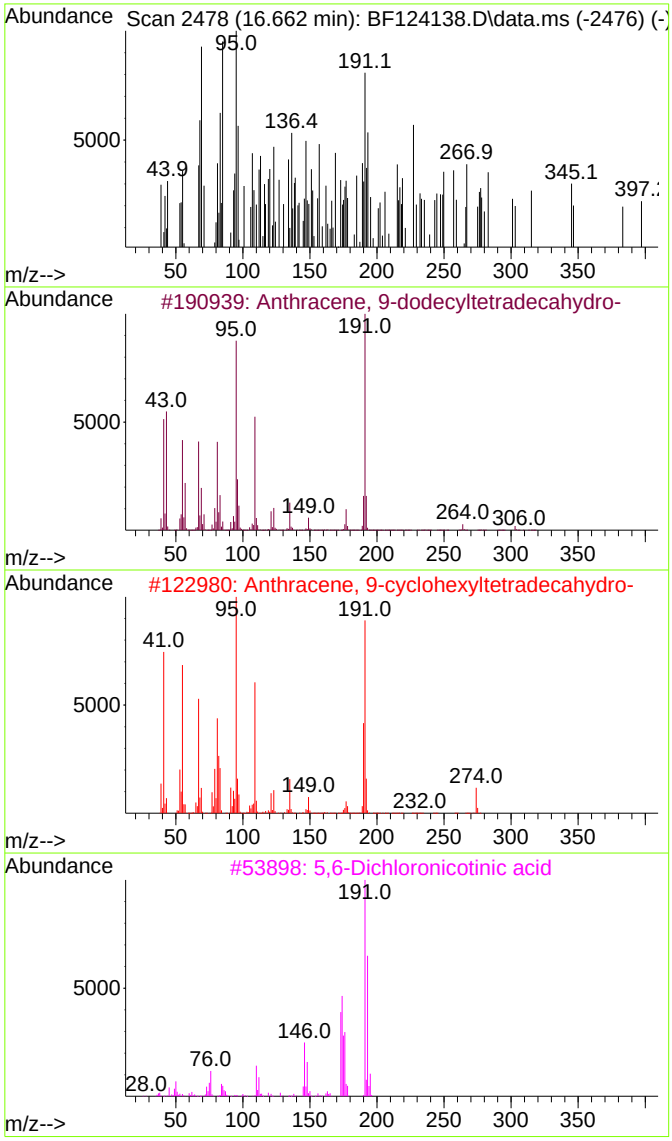
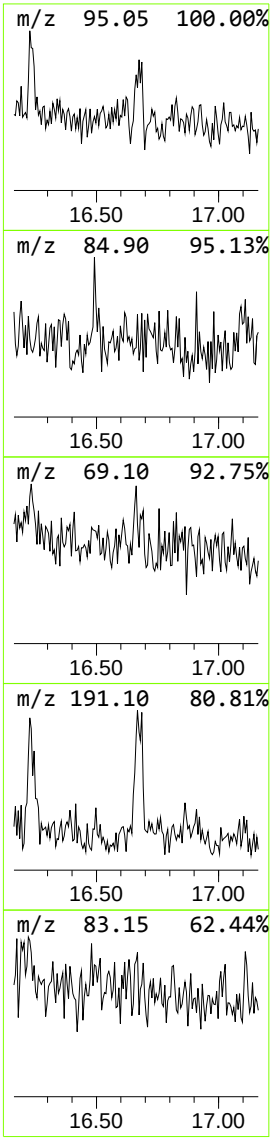
Peak Number 12 unknown16.662 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.662	4.80 ng	60327	Perylene-d12	15.515

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	16
2	Anthracene, 9-cyclohexyltetradec...	274	C20H34	055255-70-4	14
3	5,6-Dichloronicotinic acid	191	C6H3Cl2NO2	041667-95-2	11
4	Pyrrolidine, 1-(9-borabicyclo[3....	191	C12H22BN	022516-41-2	11
5	2,3-Dimethyl-7,7-tetramethylene-...	232	C13H16N2O2	1000287-32-6	10





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Operator : JU/CG
Sample : M2587-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : FK-GF-02-073-A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.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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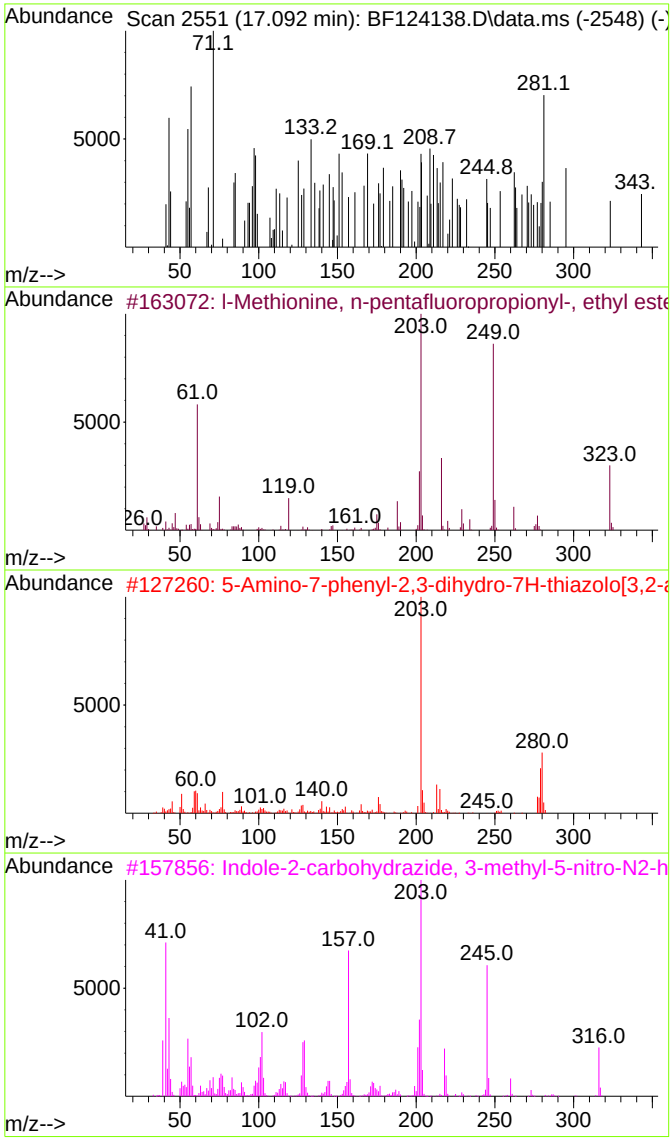
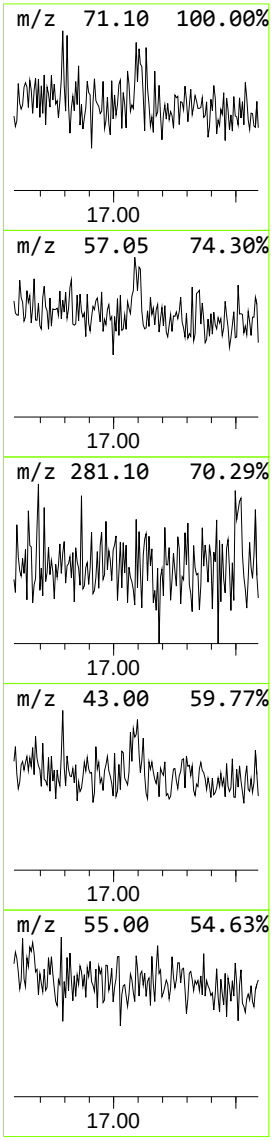
Peak Number 13 unknown17.092 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.092	2.29 ng	28789	Perylene-d12	15.515

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	l-Methionine, n-pentafluoropropi...	323	C10H14F5NO3S	1000320-90-9	22
2	5-Amino-7-phenyl-2,3-dihydro-7H-...	280	C15H12N4S	1000295-39-0	16
3	Indole-2-carbohydrazide, 3-methy...	316	C16H20N4O3	282103-55-3	12
4	5-(2-Hydroxy-4-methoxy-benzylide...	262	C12H10N2O5	1000318-35-4	10
5	4-Hexylphenyl trifluoroacetate	274	C14H17F3O2	1000373-44-0	10





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060421\
Data File : BF124138.D
Acq On : 4 Jun 2021 14:16
Operator : JU/CG
Sample : M2587-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : FK-GF-02-073-A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
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
Butane, 2-metho...	2.152	42.7	ng	637749	1	6.863	298480	20.0
Propane, 1,1-di...	2.257	2.5	ng	36530	1	6.863	298480	20.0
2-Pentanone, 4-...	5.093	2.6	ng	38994	1	6.863	298480	20.0
Ethanol, 2-(2-b...	8.045	3.6	ng	66663	2	8.145	366275	20.0
Hexadecanoic ac...	11.780	4.0	ng	66691	4	11.386	330440	20.0
unknown11.916	11.916	3.5	ng	57254	4	11.386	330440	20.0
9,12-Octadecadi...	12.469	19.4	ng	320347	4	11.386	330440	20.0
9-Octadecenoic ...	12.486	9.8	ng	162371	4	11.386	330440	20.0
unknown13.880	13.880	4.8	ng	94980	5	14.033	392642	20.0
Benzo[e]pyrene	15.380	5.7	ng	71282	6	15.515	251278	20.0
Anthracene, 9-c...	16.227	4.4	ng	54736	6	15.515	251278	20.0
unknown16.662	16.662	4.8	ng	60327	6	15.515	251278	20.0
unknown17.092	17.092	2.3	ng	28789	6	15.515	251278	20.0