LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110849.D
Acq On : 19 Nov 2018 13:31
Operator : JU/SJ
Sample : J5977-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MMTW-02(10-12)

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-BF110818.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.234	21	25	32	rVB	75111	106638	7.85%	0.726%
2	5.175	521	525	534	rBV	27583	43572	3.21%	0.297%
3	5.557	587	590	606	rBV	873501	1015166	74.74%	6.915%
4	6.563	758	761	778	rBV	1008720	1148664	84.57%	7.824%
5	6.710	782	786	793	rBV	1328410	1125395	82.85%	7.666%
6	6.939	821	825	832	rBV	414499	352664	25.96%	2.402%
7	7.092	848	851	863	rBV	988078	882538	64.97%	6.012%
8	7.504	917	921	938	rBV	677823	703728	51.81%	4.794%
9	8.222	1039	1043	1058	rBV	442508	478582	35.23%	3.260%
10	9.292	1221	1225	1241	rBV	1498869	1358285	100.00%	9.252%
11	9.692	1289	1293	1302	rBV	110307	122115	8.99%	0.832%
12	9.980	1338	1342	1353	rBV	571558	534612	39.36%	3.642%
13	10.775	1473	1477	1490	rBV	779541	761237	56.04%	5.185%
14	11.474	1592	1596	1599	rBV	547734	522278	38.45%	3.558%
15	11.498	1599	1600	1607	rVV	199329	150073	11.05%	1.022%
16	11.557	1607	1610	1618	rVB	43913	45548	3.35%	0.310%
17	11.974	1678	1681	1685	rBV	119829	125987	9.28%	0.858%
18	12.010	1685	1687	1690	rVB2	38466	33973	2.50%	0.231%
19	12.092	1698	1701	1704	rBV2	36252	40483	2.98%	0.276%
20	12.274	1728	1732	1737	rBV	17122	21631	1.59%	0.147%
21	12.527	1772	1775	1779	rBV2	28635	30173	2.22%	0.206%
22	12.692	1800	1803	1807	rBV	286289	269608	19.85%	1.837%
23	12.763	1812	1815	1819	rBV2	40196	41520	3.06%	0.283%
24	12.851	1826	1830	1831	rBV3	13957	15645	1.15%	0.107%
25	12.863	1831	1832	1836	rVB2	19201	16523	1.22%	0.113%
26	12.904	1836	1839	1840	rBV	70826	57737	4.25%	0.393%
27	12.927	1840	1843	1848	rVV	251716	240745	17.72%	1.640%
28	12.974	1848	1851	1855	rVB2	16760	17270	1.27%	0.118%
29	13.057	1861	1865	1868	rBV	1207708	1044317	76.88%	7.114%
30	13.139	1875	1879	1884	rVB6	16327	30063	2.21%	0.205%
31	13.257	1890	1899	1902	rBV3	59189	84662	6.23%	0.577%
32	13.321	1907	1910	1912	rBV2	30301	26442	1.95%	0.180%
33	13.357	1912	1916	1920	rBV2	21347	26703	1.97%	0.182%
34	13.521	1939	1944	1955	rVB	384652	409701	30.16%	2.791%

LSC Area Percent Report

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Data File : BF110849.D
Acq On : 19 Nov 2018 13:31
Operator : JU/SJ
Sample : J5977-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MMTW-02(10-12)

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-BF110818.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.604	1955	1958	1960	rBV	32917	30412	2.24%	0.207%
36	13.774	1984	1987	1990	rBV	15245	16620	1.22%	0.113%
37	13.880	2002	2005	2007	rBV	21302	17794	1.31%	0.121%
38	13.933	2011	2014	2020	rVB3	110285	122005	8.98%	0.831%
39	14.063	2028	2036	2039	rBV5	28585	72688	5.35%	0.495%
40	14.127	2042	2047	2050	rVV2	468863	577414	42.51%	3.933%
41	14.151	2050	2051	2057	rVB	133167	110535	8.14%	0.753%
42	14.245	2062	2067	2070	rVV2	62614	89794	6.61%	0.612%
43	14.292	2072	2075	2080	rVB7	16886	29835	2.20%	0.203%
44	14.480	2104	2107	2108	rBV3	14165	15695	1.16%	0.107%
45	14.515	2110	2113	2116	rVV	39928	47645	3.51%	0.325%
46	14.551	2116	2119	2123	rVB	77367	70454	5.19%	0.480%
47	14.868	2168	2173	2177	rBV	84857	100808	7.42%	0.687%
48	15.033	2198	2201	2204	rVB4	16377	21508	1.58%	0.147%
49	15.215	2225	2232	2238	rBV2	158521	365525	26.91%	2.490%
50	15.315	2246	2249	2253	rVB	27589	33540	2.47%	0.228%
51	15.521	2281	2284	2292	rBV	64564	84705	6.24%	0.577%
52	15.592	2292	2296	2303	rBV2	126899	256706	18.90%	1.749%
53	15.657	2303	2307	2311	rBV	274517	338322	24.91%	2.305%
54	15.827	2334	2336	2342	rVB6	11326	14809	1.09%	0.101%
55	16.062	2371	2376	2381	rBV	64715	103679	7.63%	0.706%
56	16.586	2461	2465	2476	rVB2	38056	71646	5.27%	0.488%
57	17.227	2567	2574	2582	rVB3	44397	111687	8.22%	0.761%
58	17.321	2584	2590	2592	rBV5	8162	16702	1.23%	0.114%
59	17.715	2651	2657	2664	rBV3	31727	75689	5.57%	0.516%

Sum of corrected areas: 14680495

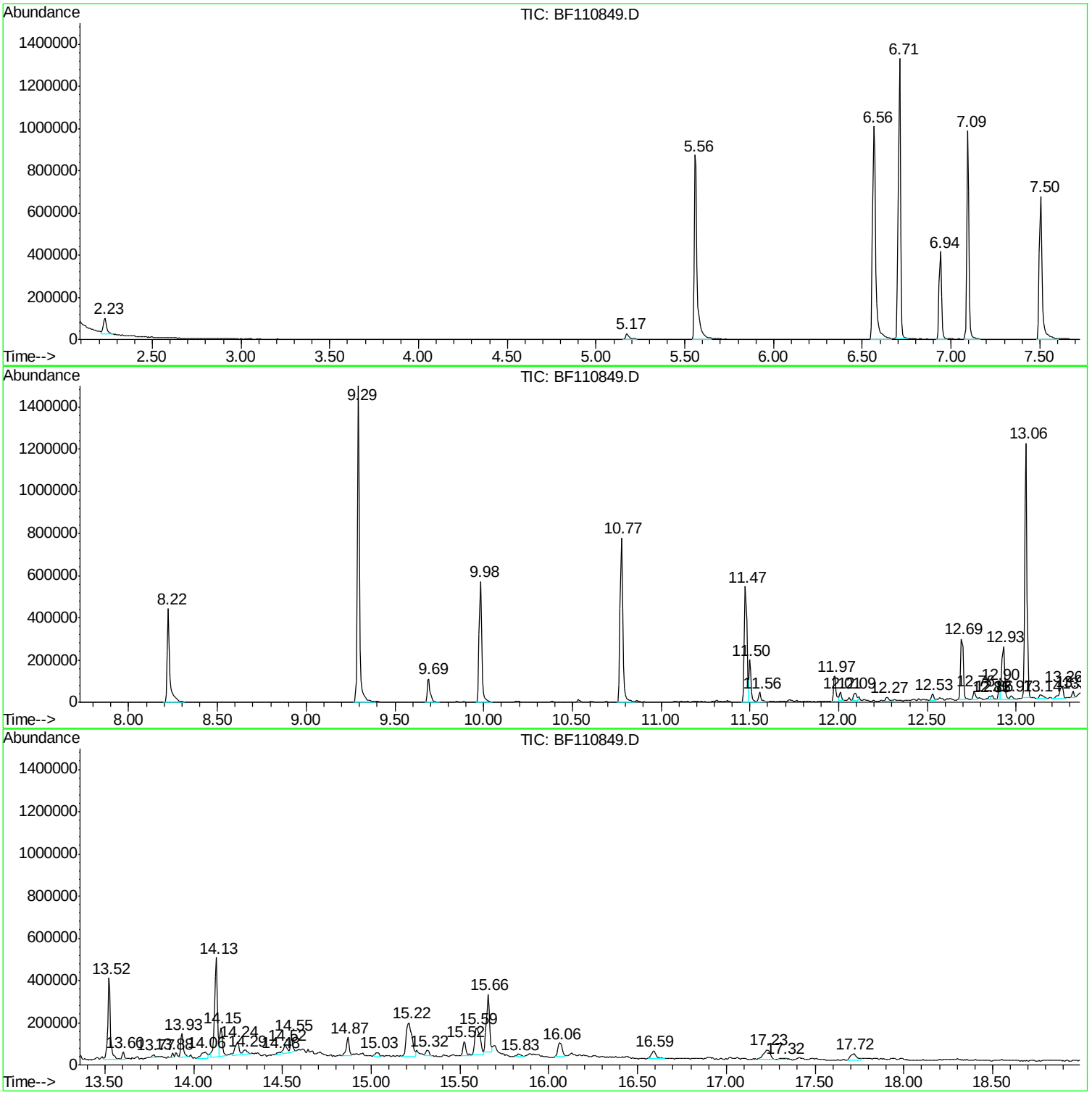
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110849.D
Acq On : 19 Nov 2018 13:31
Operator : JU/SJ
Sample : J5977-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MMTW-02(10-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110849.D
Acq On : 19 Nov 2018 13:31
Operator : JU/SJ
Sample : J5977-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MMTW-02(10-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

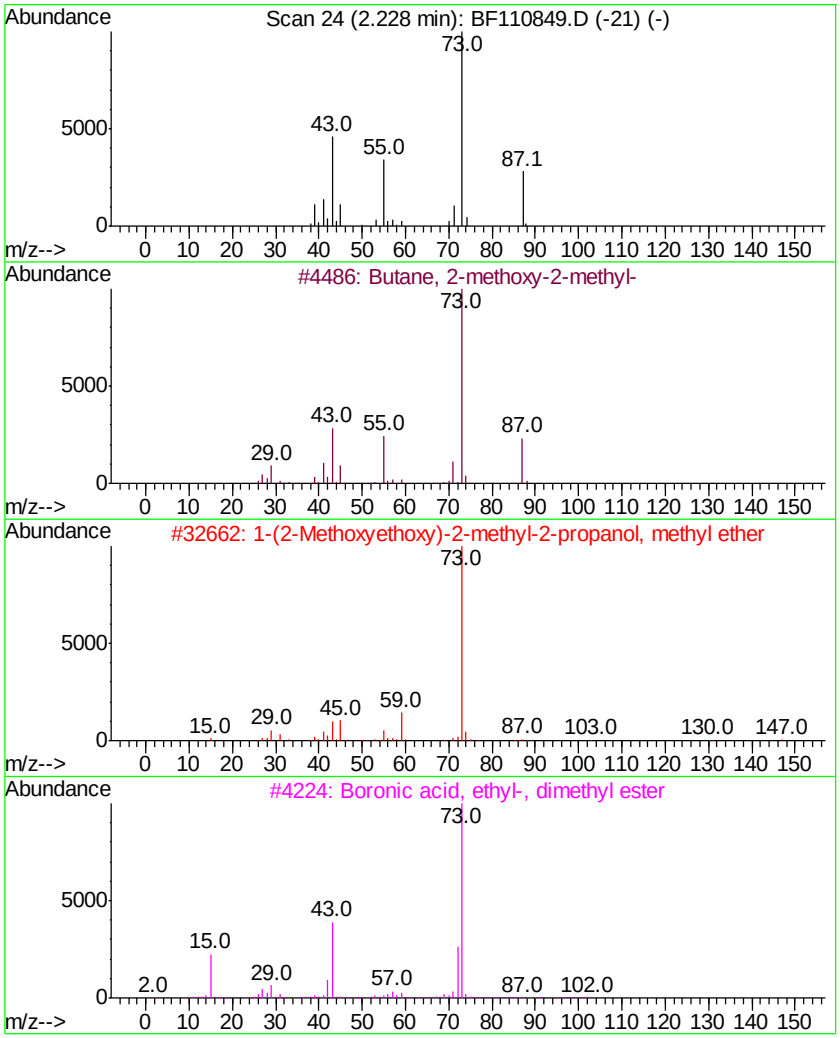
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

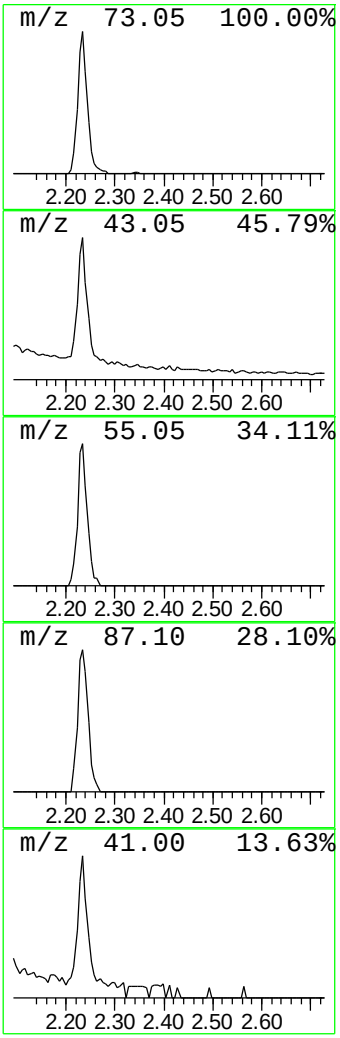
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.23	6.05 ng	106638	1,4-Dichlorobenzene-d4	6.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2	1-(2-Methoxyethoxy)-2-methyl-2-p...	162	C8H18O3	1000364-24-4	23
3	Boronic acid, ethyl-, dimethyl e...	102	C4H11BO2	007318-82-3	17
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	12
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110849.D
Acq On : 19 Nov 2018 13:31
Operator : JU/SJ
Sample : J5977-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MMTW-02(10-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

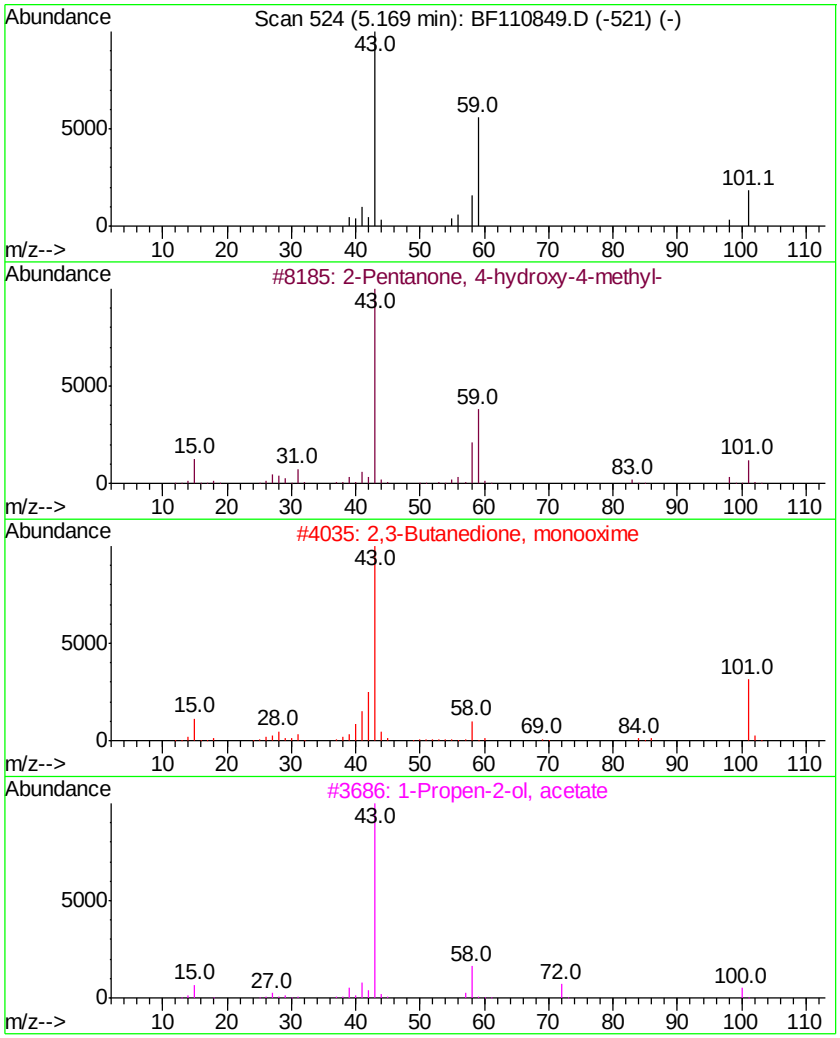
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

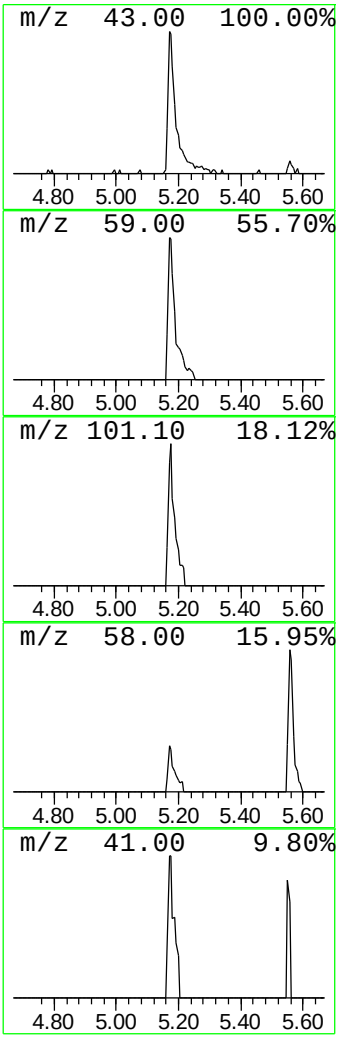
****

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	2.47 ng	43572	1,4-Dichlorobenzene-d4	6.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
5	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110849.D
Acq On : 19 Nov 2018 13:31
Operator : JU/SJ
Sample : J5977-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MMTW-02(10-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

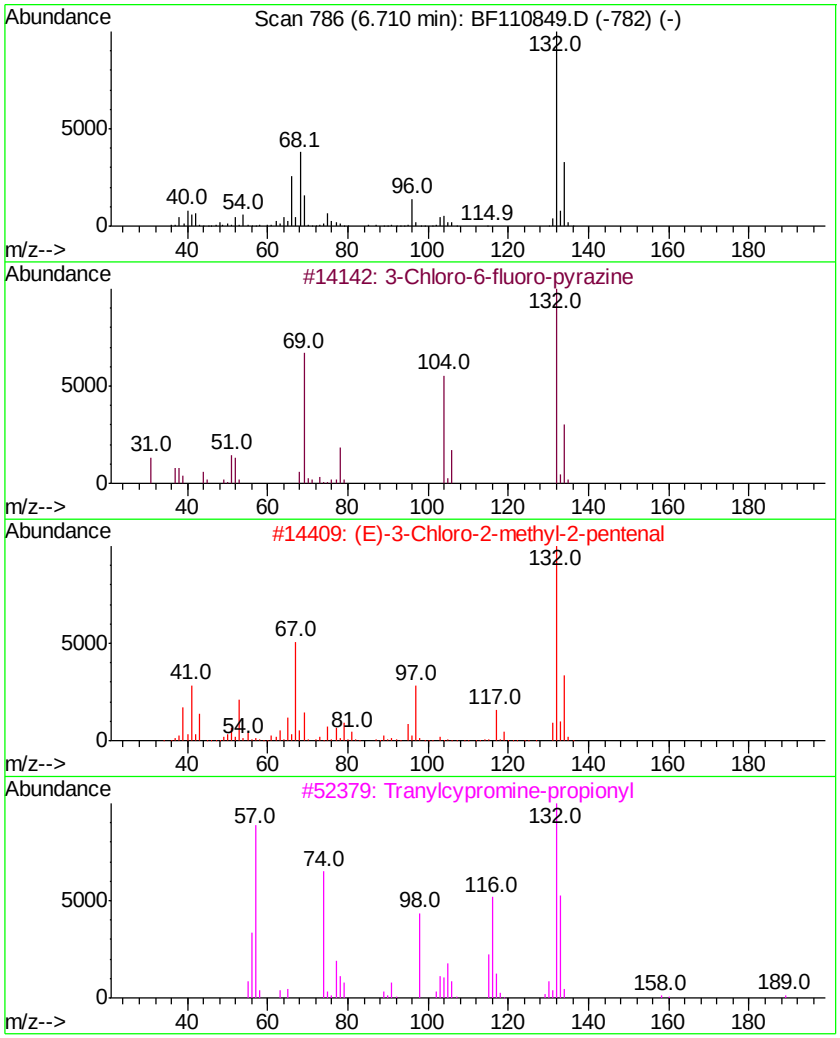
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

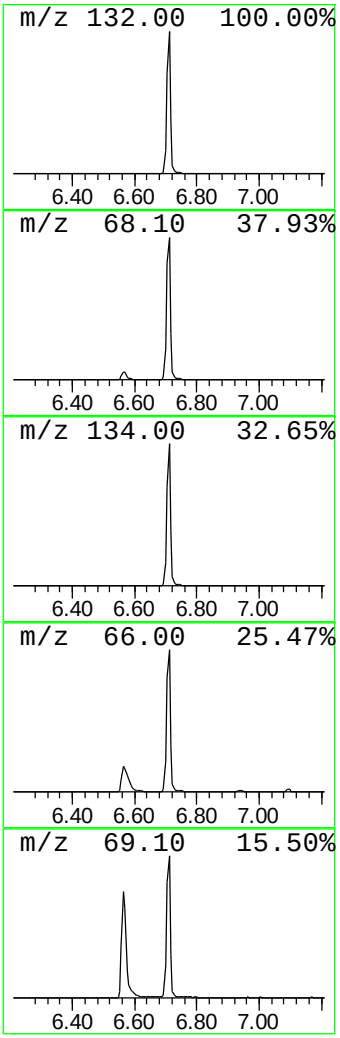
****

Peak Number 3 3-Chloro-6-fluoro-pyrazine Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.71	63.82 ng	1125400	1,4-Dichlorobenzene-d4	6.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	16
4	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5	4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110849.D
Acq On : 19 Nov 2018 13:31
Operator : JU/SJ
Sample : J5977-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MMTW-02(10-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

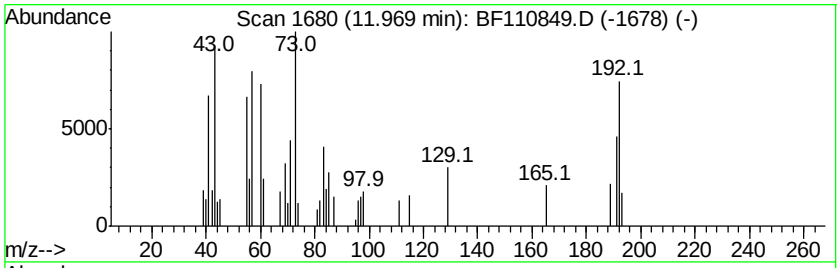
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

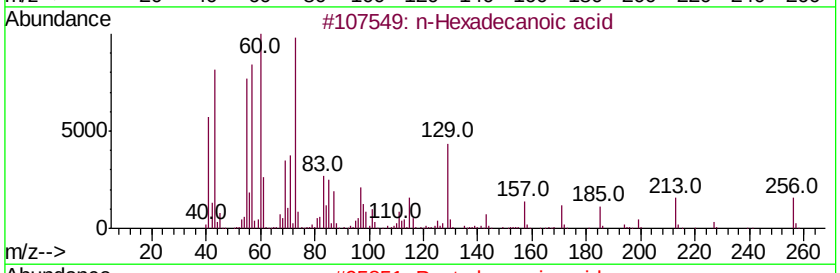
****

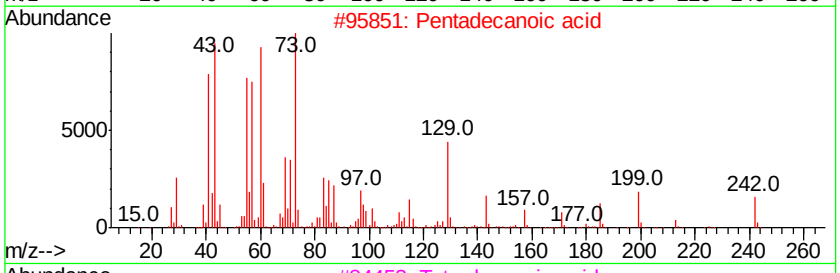
Peak Number 5 n-Hexadecanoic acid Concentration Rank 7

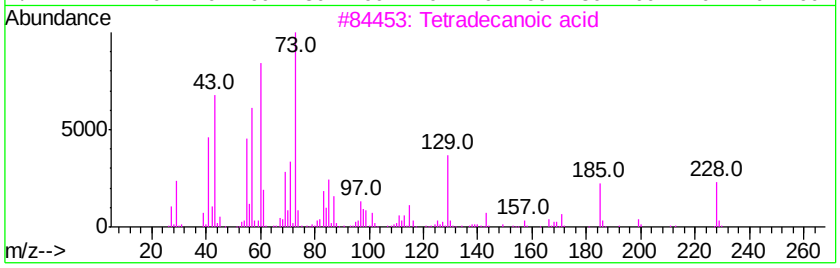
R.T.	EstConc	Area	Relative to ISTD	R.T.
11.97	4.82 ng	125987	Phenanthrene-d10	11.47

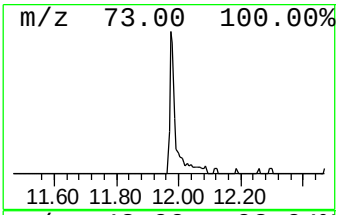
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	64
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	52
4	Tridecanoic acid	214	C13H26O2	000638-53-9	49
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	35

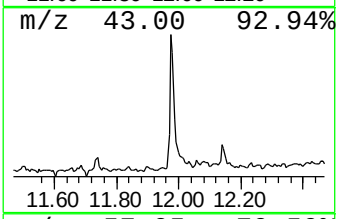


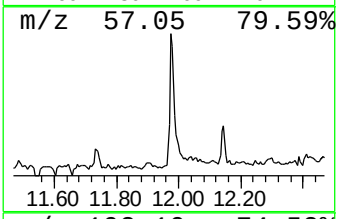


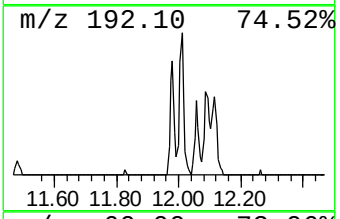


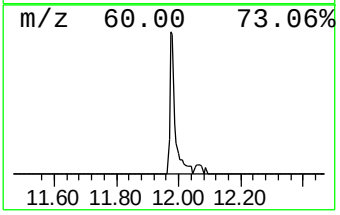












Library Search Compound Report

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Data File : BF110849.D
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Instrument : BNA_F
ClientSampleId : MMTW-02(10-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

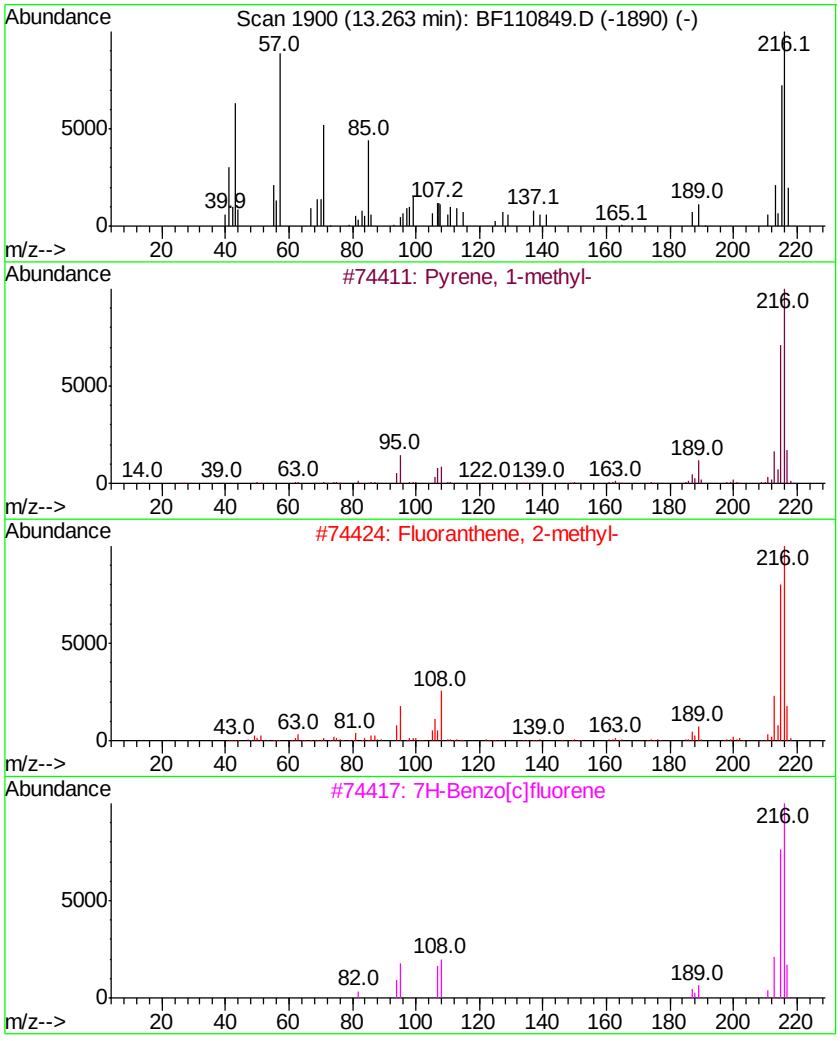
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

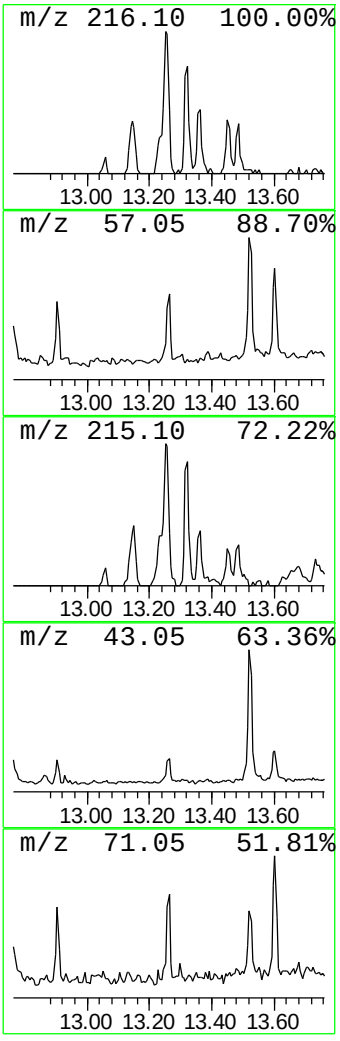
****

Peak Number 6 Pyrene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.26	2.93 ng	84662	Chrysene-d12	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	81
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	76
3	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	70
5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	68





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110849.D
Acq On : 19 Nov 2018 13:31
Operator : JU/SJ
Sample : J5977-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MMTW-02(10-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

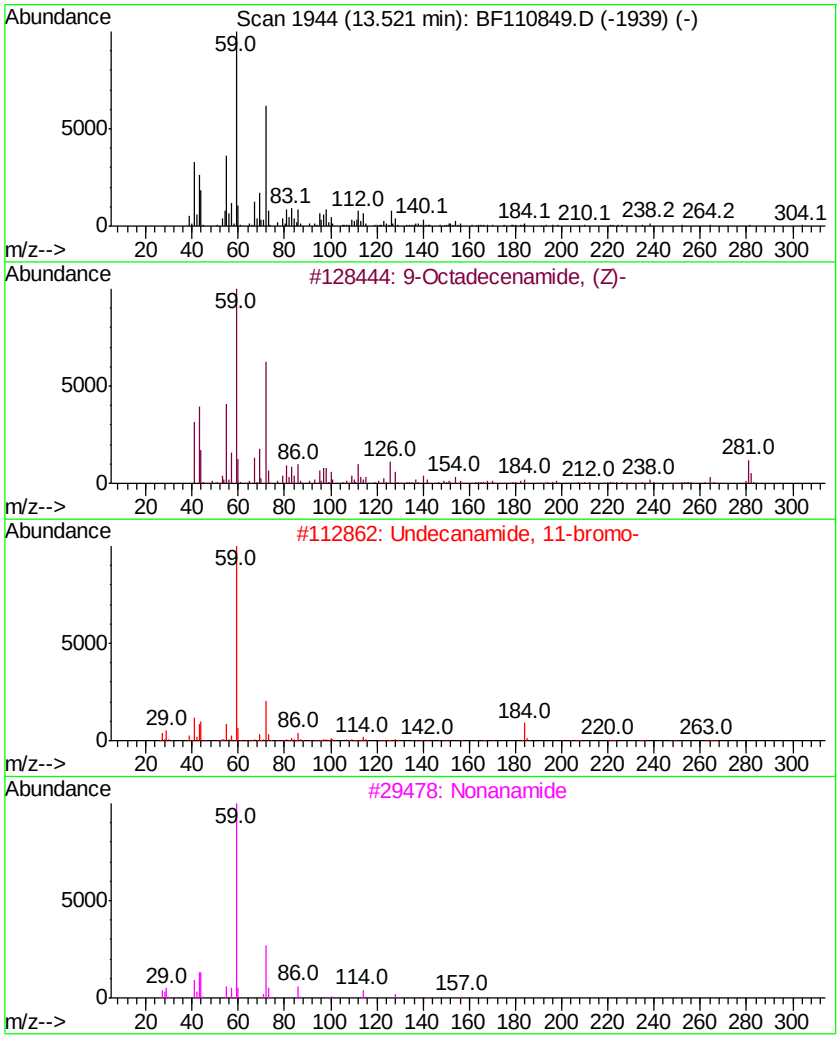
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

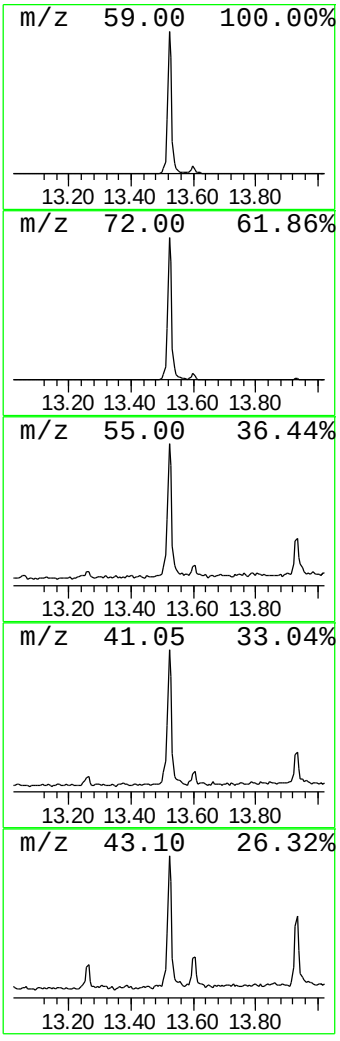
****

Peak Number 7 9-Octadecenamide, (Z)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.52	14.19 ng	409701	Chrysene-d12	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	96
2	Undecanamide, 11-bromo-	263	C11H22BrNO	005875-26-3	59
3	Nonanamide	157	C9H19NO	001120-07-6	59
4	Nonadecanamide	297	C19H39NO	058185-32-3	56
5	Octanamide	143	C8H17NO	000629-01-6	53





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
Data File : BF110849.D
Acq On : 19 Nov 2018 13:31
Operator : JU/SJ
Sample : J5977-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MMTW-02(10-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

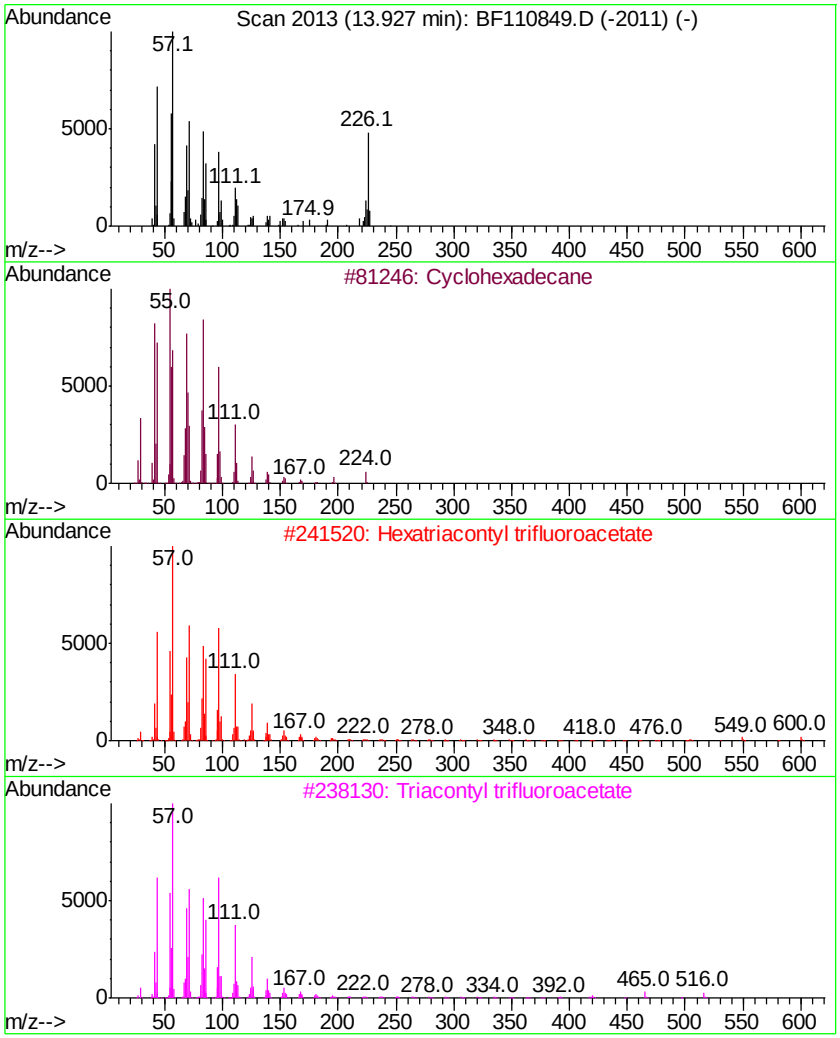
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

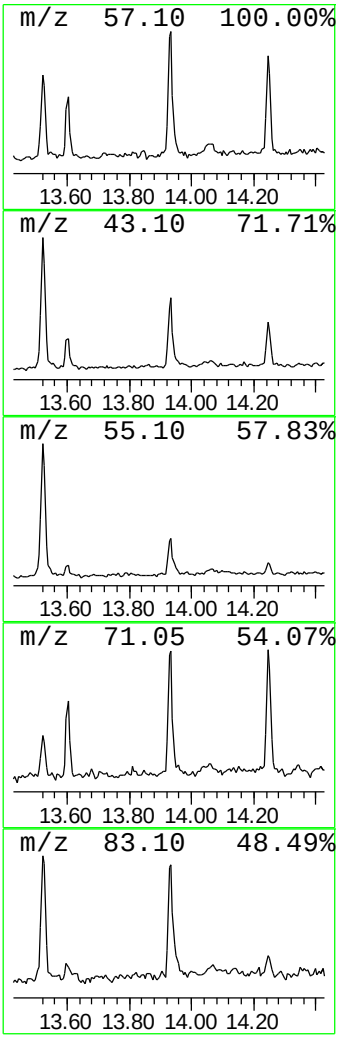
****

Peak Number 8 Cyclohexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.93	4.23 ng	122005	Chrysene-d12	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	86
2	Hexatriacontyl trifluoroacetate	619	C38H73F3O2	1000351-88-6	62
3	Triacontyl trifluoroacetate	534	C32H61F3O2	1000351-75-7	62
4	Octacosyl heptafluorobutyrate	606	C32H57F7O2	1000351-83-6	62
5	1-Decanol, 2-octyl-	270	C18H38O	045235-48-1	62





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Sample : J5977-11
Misc :
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Instrument : BNA_F
ClientSampleId : MMTW-02(10-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M
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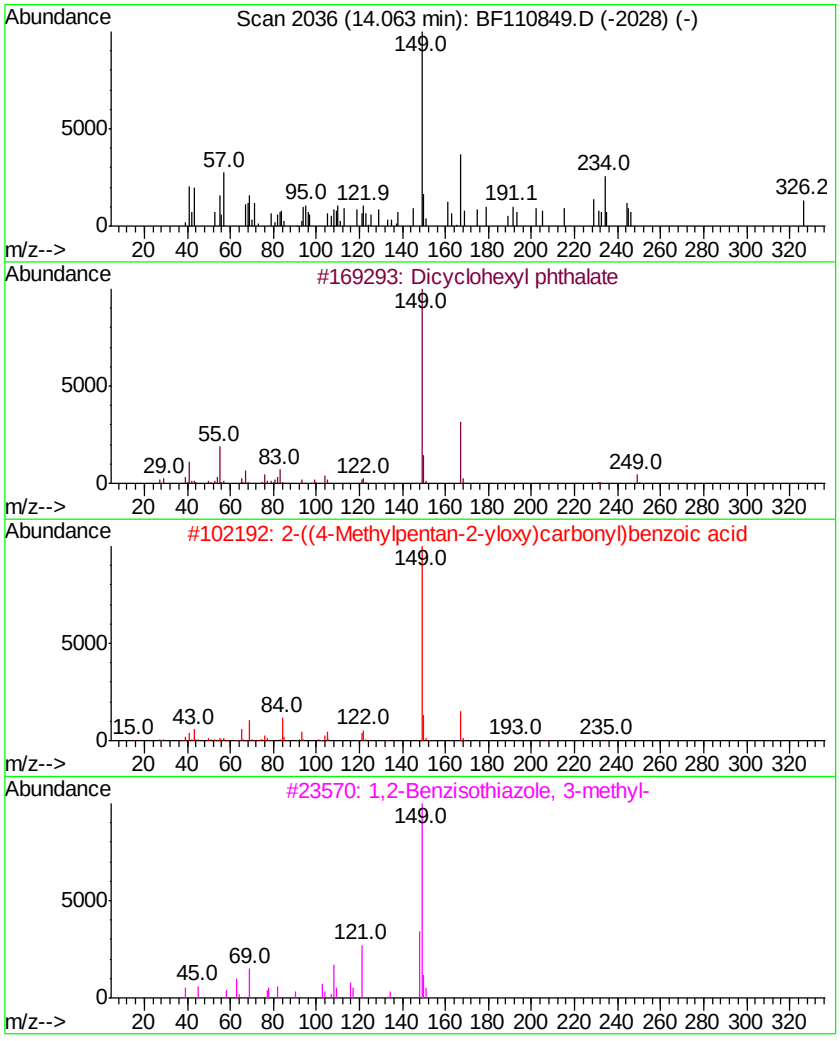
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

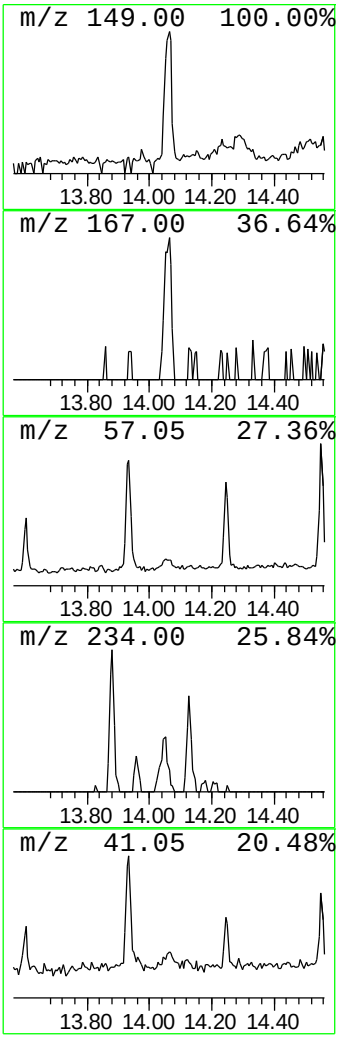
****

Peak Number 9 unknown14.06 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.06	2.52 ng	72688	Chrysene-d12	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dicyclohexyl phthalate	330	C20H26O4	000084-61-7	43
2	2-((4-Methylpentan-2-yloxy)carbo...	250	C14H18O4	1000373-82-5	43
3	1,2-Benzisothiazole, 3-methyl-	149	C8H7NS	006187-89-9	38
4	2-(Decyloxycarbonyl)benzoic acid	306	C18H26O4	1000373-53-2	38
5	Phthalic acid, undec-2-en-1-yl u...	472	C30H48O4	1000315-42-2	38





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Data File : BF110849.D
Acq On : 19 Nov 2018 13:31
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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MMTW-02(10-12)

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TIC Library : C:\DATABASE\NIST11.L
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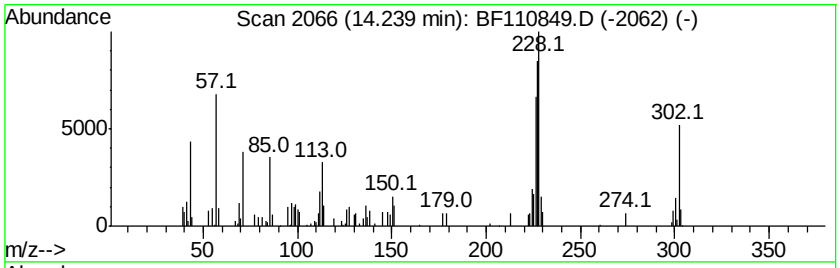
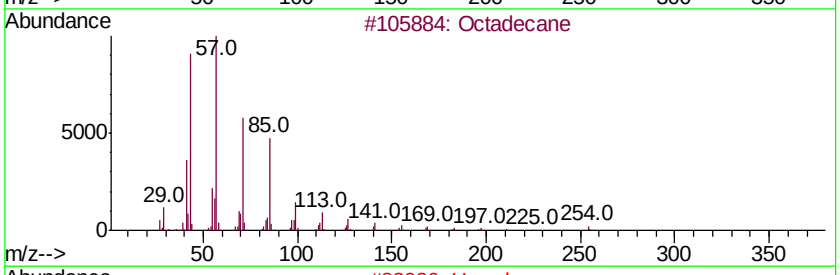
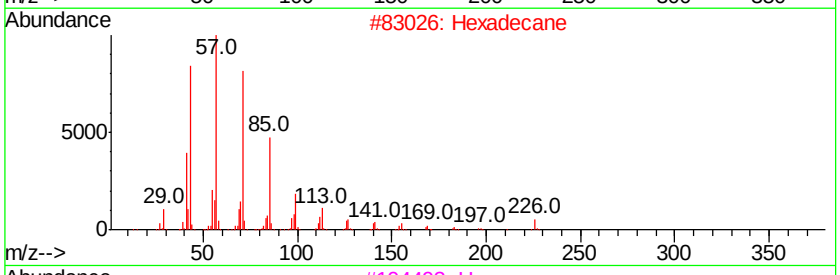
Peak Number 10 Octadecane Concentration Rank 11

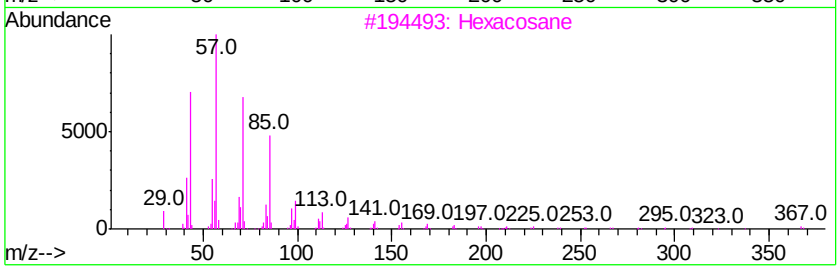
R.T.	EstConc	Area	Relative to ISTD	R.T.
14.24	3.11 ng	89794	Chrysene-d12	14.13

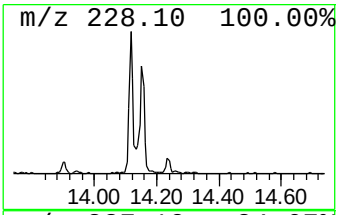
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	90
2	Hexadecane	226	C16H34	000544-76-3	87
3	Hexacosane	366	C26H54	000630-01-3	49
4	Hentriacontane	437	C31H64	000630-04-6	49
5	Heptadecane	240	C17H36	000629-78-7	46

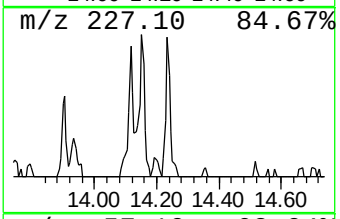


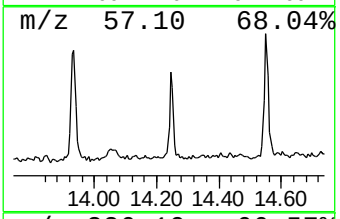


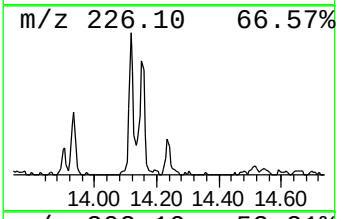


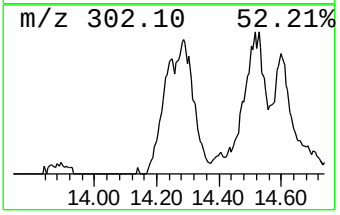












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Sample : J5977-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MMTW-02(10-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

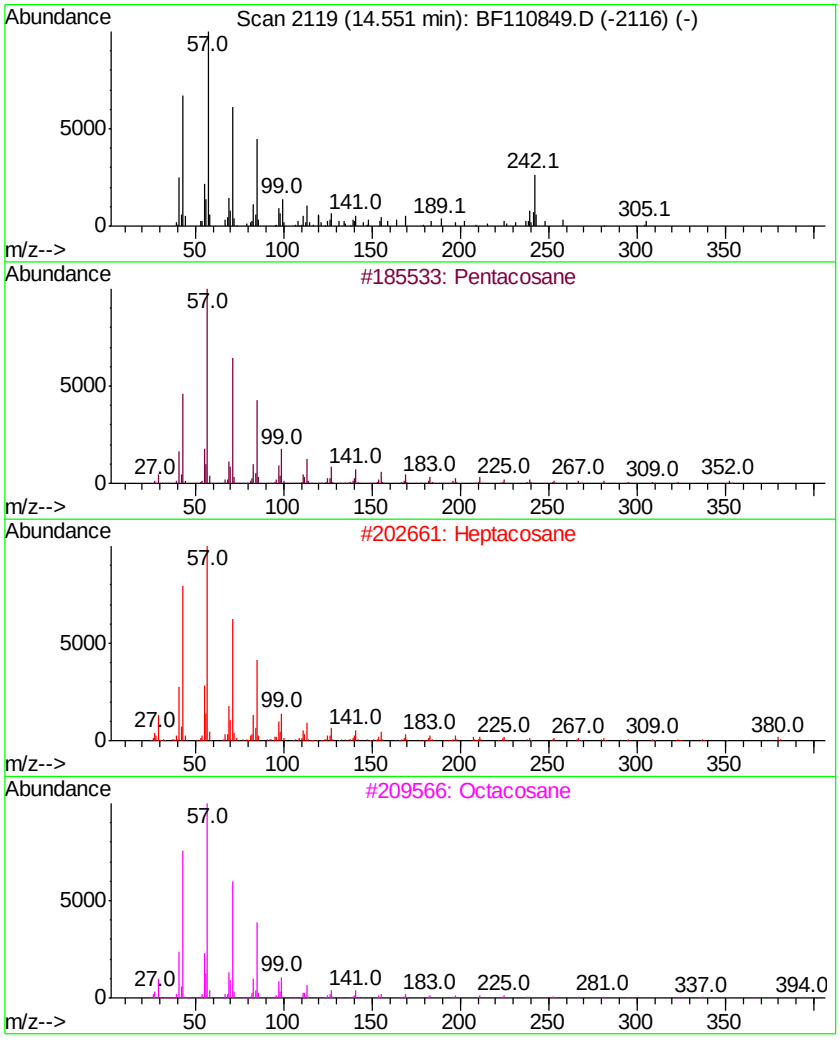
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

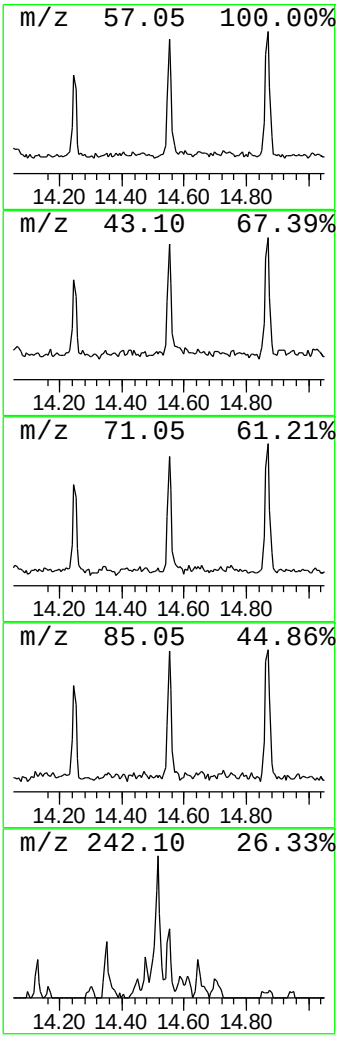
****

Peak Number 11 Hexatriacontane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.55	2.44 ng	70454	Chrysene-d12	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane	352	C25H52	000629-99-2	81
2	Heptacosane	380	C27H56	000593-49-7	81
3	Octacosane	394	C28H58	000630-02-4	81
4	Hexatriacontane	507	C36H74	000630-06-8	81
5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	81





Library Search Compound Report

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Instrument : BNA_F
ClientSampleId : MMTW-02(10-12)

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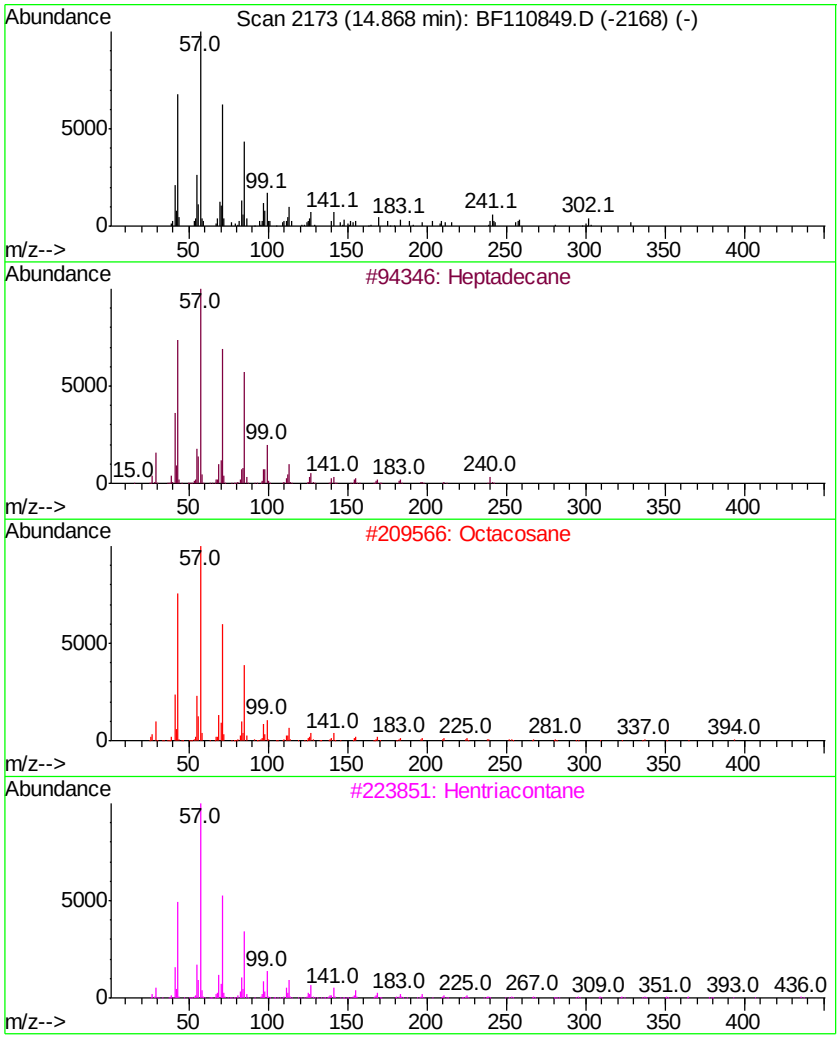
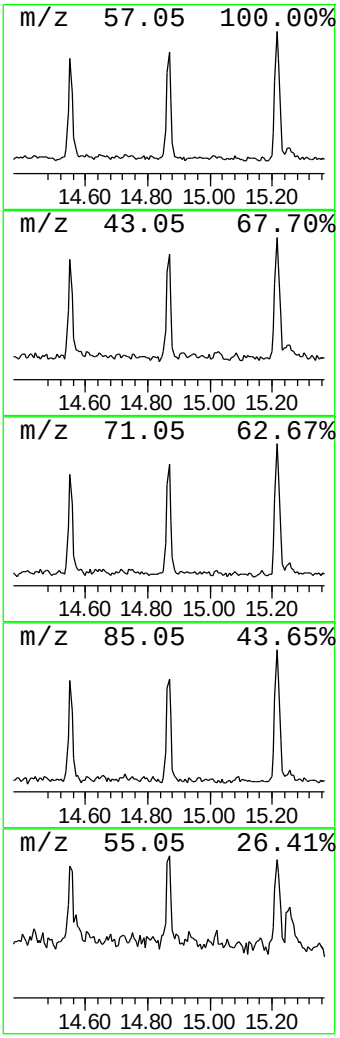
Peak Number 12 Heptadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	3.49 ng	100808	Chrysene-d12	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	97
2	Octacosane	394	C28H58	000630-02-4	87
3	Hentriacontane	437	C31H64	000630-04-6	87
4	Tetracosane	338	C24H50	000646-31-1	87
5	Hexacosane	366	C26H54	000630-01-3	86





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Sample : J5977-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MMTW-02(10-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M
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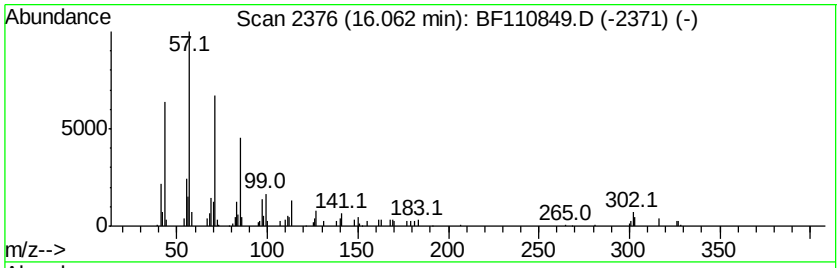
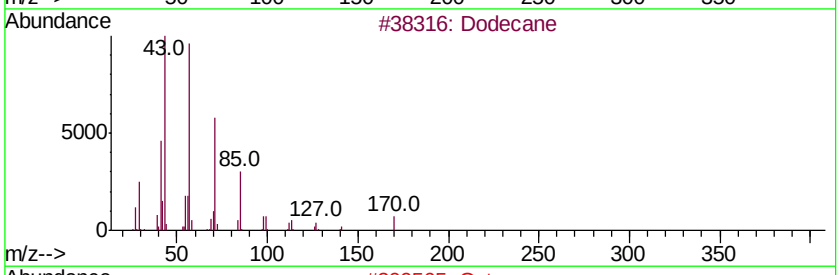
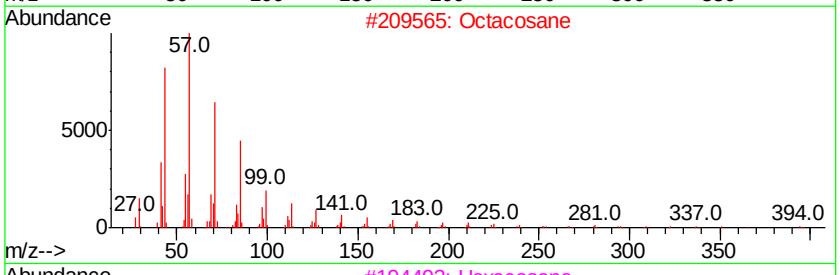
Peak Number 14 Dodecane Concentration Rank 4

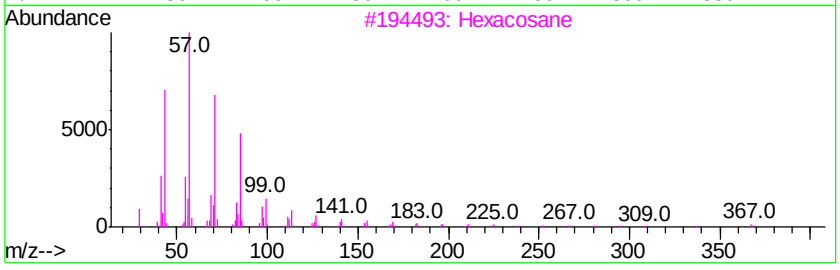
R.T.	EstConc	Area	Relative to ISTD	R.T.
16.06	6.13 ng	103679	Perylene-d12	15.66

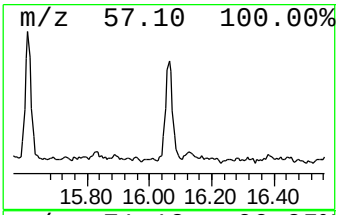
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane	170	C12H26	000112-40-3	89
2	Octacosane	394	C28H58	000630-02-4	87
3	Hexacosane	366	C26H54	000630-01-3	87
4	Heneicosane	296	C21H44	000629-94-7	87
5	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	87

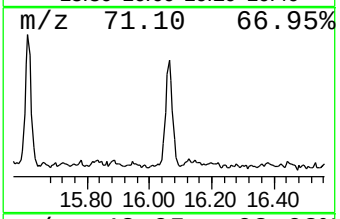


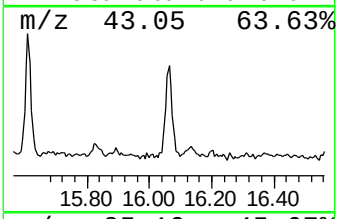


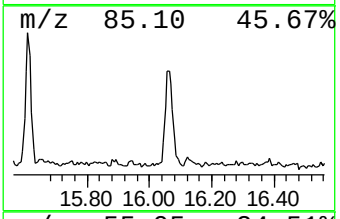


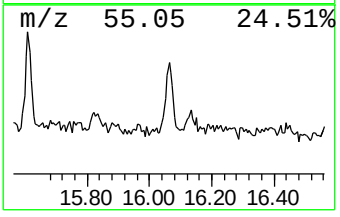












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TIC Library : C:\DATABASE\NIST11.L
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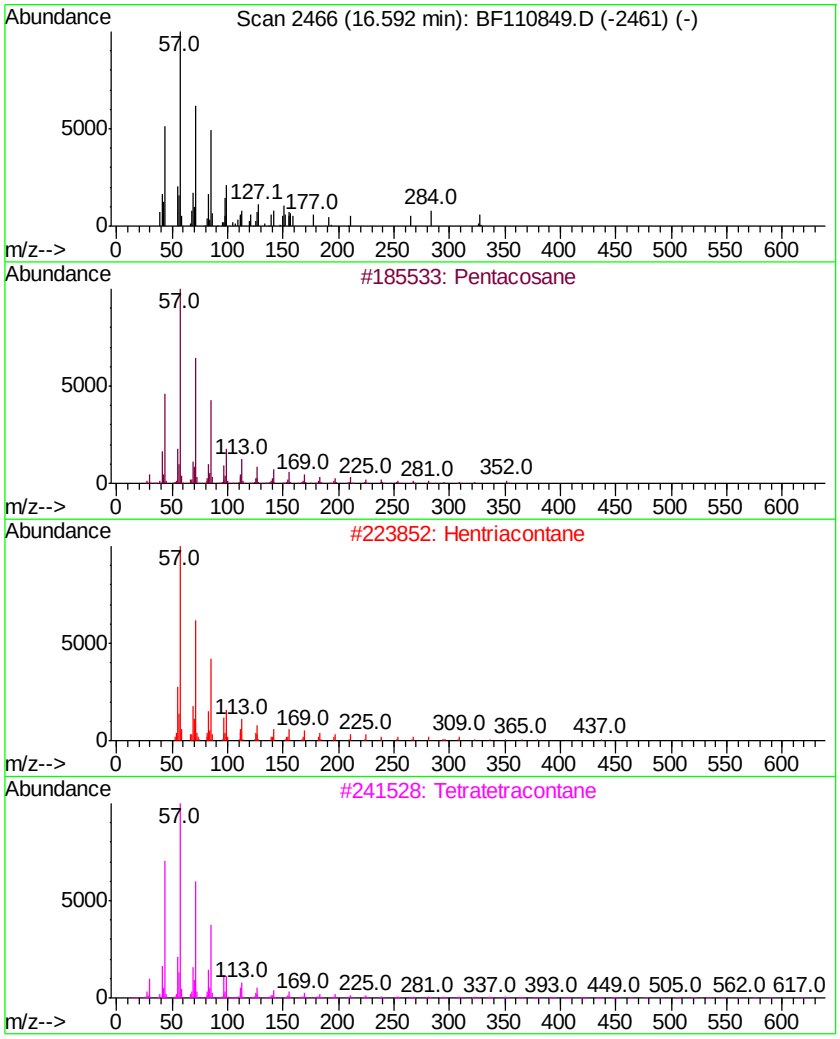
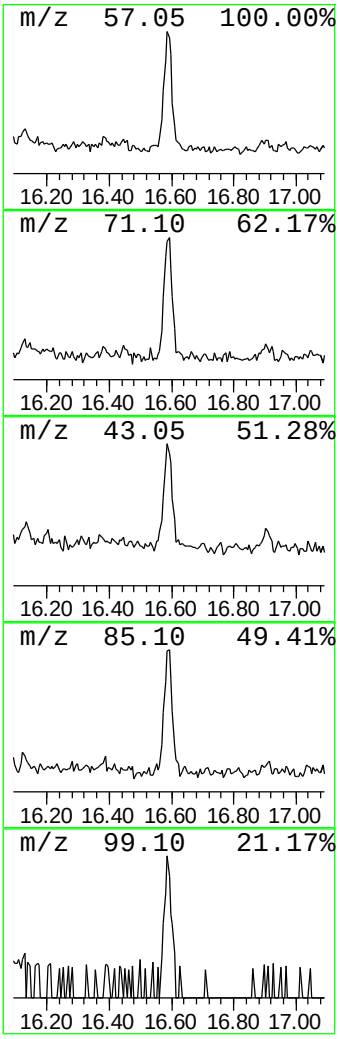
****

Peak Number 15 Pentacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.59	4.24 ng	71646	Perylene-d12	15.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane	352	C25H52	000629-99-2	72
2	Hentriacontane	437	C31H64	000630-04-6	72
3	Tetratetracontane	619	C44H90	007098-22-8	64
4	Docosane	310	C22H46	000629-97-0	64
5	Sulfurous acid, dodecyl 2-propyl...	292	C15H32O3S	1000309-12-3	64





Tentatively Identified Compound (LSC) summary

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TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methoxy...	2.23	6.0	ng	106638	1	6.94	352664	20.0
2-Pentanone, 4-hy...	5.17	2.5	ng	43572	1	6.94	352664	20.0
3-Chloro-6-fluoro...	6.71	63.8	ng	1125400	1	6.94	352664	20.0
n-Hexadecanoic acid	11.97	4.8	ng	125987	4	11.47	522278	20.0
Pyrene, 1-methyl-	13.26	2.9	ng	84662	5	14.13	577414	20.0
9-Octadecenamide,...	13.52	14.2	ng	409701	5	14.13	577414	20.0
Cyclohexadecane	13.93	4.2	ng	122005	5	14.13	577414	20.0
unknown14.06	14.06	2.5	ng	72688	5	14.13	577414	20.0
Octadecane	14.24	3.1	ng	89794	5	14.13	577414	20.0
Hexatriacontane	14.55	2.4	ng	70454	5	14.13	577414	20.0
Heptadecane	14.87	3.5	ng	100808	5	14.13	577414	20.0
Dodecane	16.06	6.1	ng	103679	6	15.66	338322	20.0
Pentacosane	16.59	4.2	ng	71646	6	15.66	338322	20.0