LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060119\
Data File : BF114744.D
Acq On : 1 Jun 2019 9:34
Operator : HP/JU
Sample : K3097-23
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 0443-PR-6(PASSAIC)

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-BF052919.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	3.675	261	270	276	rVB2	46552	96757	1.04%	0.093%
2	5.310	542	548	551	rBV	7811733	8672897	93.41%	8.348%
3	6.334	715	722	725	rBV	7226000	9143705	98.48%	8.801%
4	6.469	739	745	748	rBV	8351600	9285040	100.00%	8.937%
5	6.692	780	783	787	rBV	2601247	2173711	23.41%	2.092%
6	6.851	806	810	813	rBV	7338334	6733032	72.51%	6.481%
7	6.875	813	814	819	rVB2	119567	121333	1.31%	0.117%
8	7.257	870	879	882	rBV	5544246	5893725	63.48%	5.673%
9	7.345	890	894	899	rVB6	69997	120223	1.29%	0.116%
10	7.622	935	941	946	rVB8	81185	136969	1.48%	0.132%
11	7.975	997	1001	1003	rBV	3323422	2988247	32.18%	2.876%
12	7.992	1003	1004	1007	rVB	856030	542086	5.84%	0.522%
13	8.057	1013	1015	1017	rVB	141219	99224	1.07%	0.096%
14	8.416	1072	1076	1085	rVB2	196347	288773	3.11%	0.278%
15	8.527	1091	1095	1101	rBV6	77774	125621	1.35%	0.121%
16	8.680	1117	1121	1125	rVB	190542	202126	2.18%	0.195%
17	8.780	1135	1138	1142	rVB	236335	208206	2.24%	0.200%
18	9.010	1174	1177	1179	rBV	163881	121025	1.30%	0.116%
19	9.051	1179	1184	1187	rBV	8868366	8405597	90.53%	8.091%
20	9.145	1194	1200	1203	rBV5	108083	187983	2.02%	0.181%
21	9.239	1214	1216	1222	rVB	80477	115558	1.24%	0.111%
22	9.304	1224	1227	1230	rBV	94974	125858	1.36%	0.121%
23	9.380	1237	1240	1243	rBV	197295	184326	1.99%	0.177%
24	9.439	1247	1250	1253	rVV	1600090	1278194	13.77%	1.230%
25	9.463	1253	1254	1256	rVV	247937	155747	1.68%	0.150%
26	9.498	1256	1260	1262	rVB2	104535	97596	1.05%	0.094%
27	9.574	1270	1273	1276	rBV	160735	160264	1.73%	0.154%
28	9.639	1281	1284	1287	rVB	160182	146083	1.57%	0.141%
29	9.722	1293	1298	1301	rBV	3485587	3306248	35.61%	3.182%
30	9.751	1301	1303	1306	rVB	312660	239386	2.58%	0.230%
31	9.822	1311	1315	1321	rVB3	67783	136872	1.47%	0.132%
32	9.922	1329	1332	1336	rVB2	182341	191797	2.07%	0.185%
33	9.963	1336	1339	1341	rVB	133024	110347	1.19%	0.106%
34	9.998	1341	1345	1347	rBV3	176779	205952	2.22%	0.198%

LSC Area Percent Report

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Data File : BF114744.D
Acq On : 1 Jun 2019 9:34
Operator : HP/JU
Sample : K3097-23
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 0443-PR-6(PASSAIC)

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-BF052919.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	10.033	1348	1351	1354	rVB3	80559	110624	1.19%	0.106%
36	10.127	1363	1367	1372	rVB5	98852	166525	1.79%	0.160%
37	10.263	1387	1390	1392	rBV	276425	238076	2.56%	0.229%
38	10.392	1407	1412	1414	rVB2	202141	248233	2.67%	0.239%
39	10.504	1426	1431	1434	rBV	5866298	5666348	61.03%	5.454%
40	10.657	1454	1457	1463	rVB2	310051	315679	3.40%	0.304%
41	10.804	1479	1482	1485	rVV3	96812	118256	1.27%	0.114%
42	10.833	1485	1487	1490	rVV3	125303	122218	1.32%	0.118%
43	10.892	1494	1497	1500	rVB3	87964	94539	1.02%	0.091%
44	11.116	1532	1535	1539	rVB	257693	242992	2.62%	0.234%
45	11.192	1544	1548	1550	rVV	3797000	3299582	35.54%	3.176%
46	11.216	1550	1552	1555	rVV	1307419	980550	10.56%	0.944%
47	11.263	1557	1560	1564	rVB	441925	409755	4.41%	0.394%
48	11.304	1564	1567	1570	rBV2	108861	130021	1.40%	0.125%
49	11.668	1626	1629	1631	rBV3	88184	108582	1.17%	0.105%
50	11.692	1631	1633	1635	rVV	312860	229943	2.48%	0.221%
51	11.739	1635	1641	1644	rVV2	2102236	2743121	29.54%	2.640%
52	11.798	1648	1651	1654	rBV	441839	454636	4.90%	0.438%
53	11.986	1680	1683	1685	rBV	146973	115824	1.25%	0.111%
54	12.239	1723	1726	1729	rBV	258672	265679	2.86%	0.256%
55	12.274	1730	1732	1734	rVB2	107212	103329	1.11%	0.099%
56	12.398	1750	1753	1756	rBV	1119224	1011691	10.90%	0.974%
57	12.445	1758	1761	1766	rVV	657078	774292	8.34%	0.745%
58	12.521	1769	1774	1786	rVV	2688194	3920634	42.23%	3.774%
59	12.621	1787	1791	1793	rVV	1096075	1049143	11.30%	1.010%
60	12.645	1793	1795	1798	rVB2	108714	106979	1.15%	0.103%
61	12.774	1813	1817	1821	rBV	7670050	7845621	84.50%	7.552%
62	12.951	1845	1847	1852	rVB	313760	285123	3.07%	0.274%
63	13.015	1856	1858	1862	rVB2	293466	260632	2.81%	0.251%
64	13.051	1862	1864	1868	rBV	174898	189527	2.04%	0.182%
65	13.174	1883	1885	1890	rVB	164268	154688	1.67%	0.149%
66	13.680	1968	1971	1977	rVB2	391027	551121	5.94%	0.530%
67	13.810	1988	1993	1996	rBV2	2620026	2982504	32.12%	2.871%
68	13.833	1996	1997	2005	rVB	428825	366618	3.95%	0.353%
69	14.004	2024	2026	2030	rVB	134309	107834	1.16%	0.104%
70	14.198	2057	2059	2061	rVB	136540	100762	1.09%	0.097%
71	14.309	2076	2078	2081	rBV3	195211	179158	1.93%	0.172%

LSC Area Percent Report

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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 0443-PR-6(PASSAIC)

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

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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052919.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	14.604	2125	2128	2134	rVB2	283889	276679	2.98%	0.266%
73	14.674	2137	2140	2143	rVB	156651	159332	1.72%	0.153%
74	14.809	2160	2163	2166	rBV	255364	365781	3.94%	0.352%
75	14.915	2177	2181	2184	rVB2	380318	442668	4.77%	0.426%
76	15.086	2207	2210	2215	rVB	190576	196925	2.12%	0.190%
77	15.139	2216	2219	2223	rVV	316205	331226	3.57%	0.319%
78	15.198	2225	2229	2233	rVV	2020820	2395573	25.80%	2.306%
79	15.262	2237	2240	2244	rVB	288628	327095	3.52%	0.315%
80	15.662	2303	2308	2312	rBV	263369	391317	4.21%	0.377%
81	16.121	2382	2386	2392	rVB	186270	318689	3.43%	0.307%
82	16.509	2449	2452	2461	rVB2	159552	290339	3.13%	0.279%
83	16.656	2474	2477	2485	rVB4	99850	183029	1.97%	0.176%
84	16.903	2515	2519	2526	rVB2	103605	164238	1.77%	0.158%

Sum of corrected areas: 103894238

LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060119\
Data File : BF114744.D
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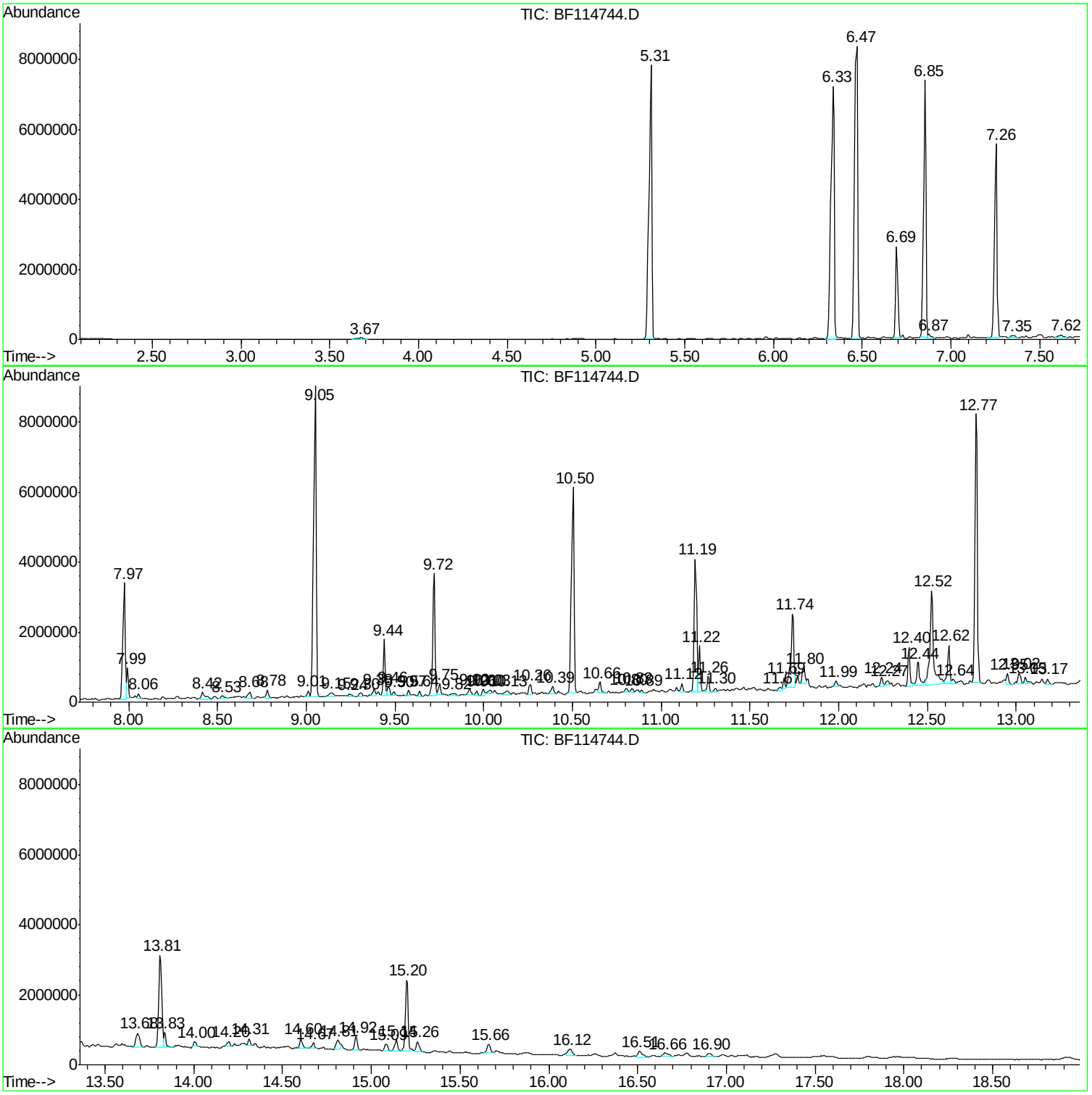
Operator : HP/JU
Sample : K3097-23
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 0443-PR-6(PASSAIC)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052919.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 F\DATA\BF060119\
Data File : BF114744.D
Acq On : 1 Jun 2019 9:34
Operator : HP/JU
Sample : K3097-23
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 0443-PR-6(PASSAIC)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

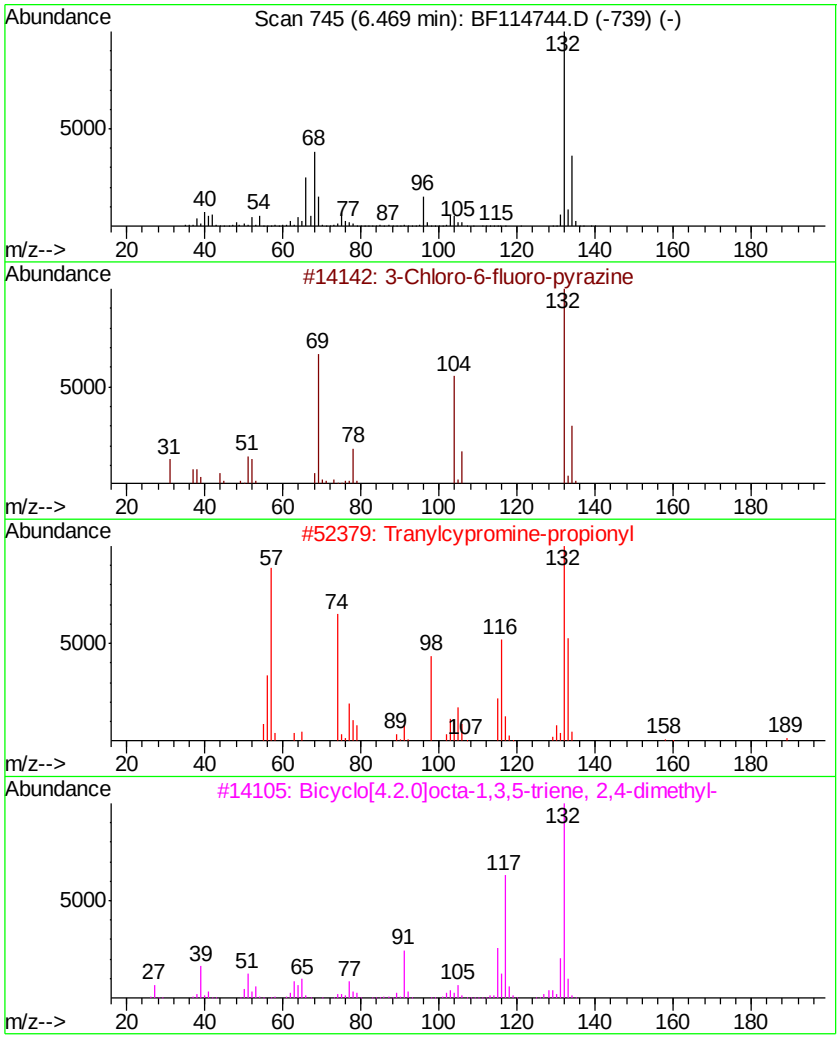
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

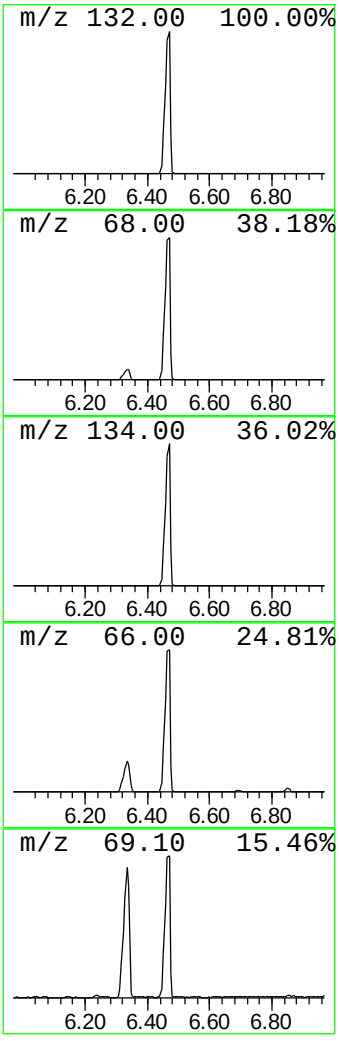
****

Peak Number 1 unknown6.47 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	85.43 ng	9285040	1,4-Dichlorobenzene-d4	6.69

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060119\
Data File : BF114744.D
Acq On : 1 Jun 2019 9:34
Operator : HP/JU
Sample : K3097-23
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 0443-PR-6(PASSAIC)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

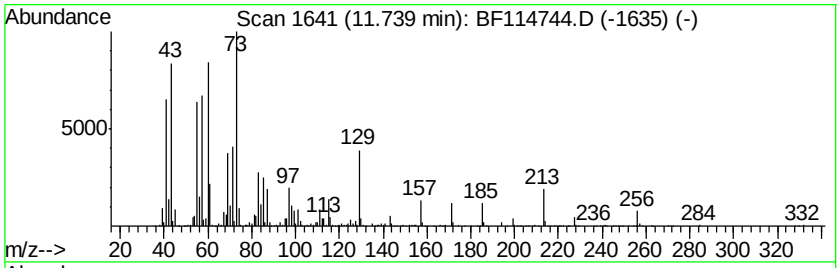
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

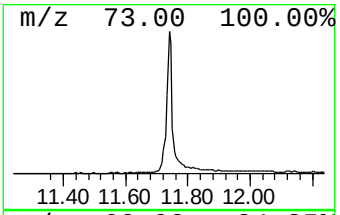
****

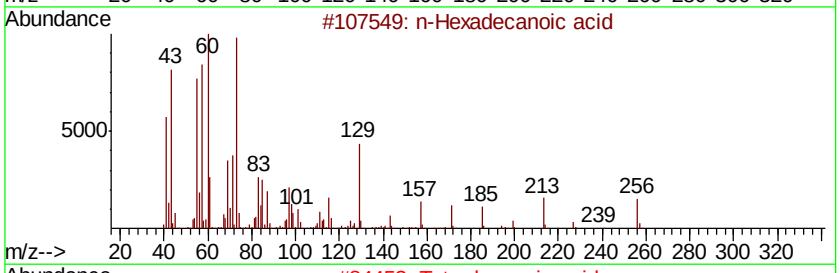
Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

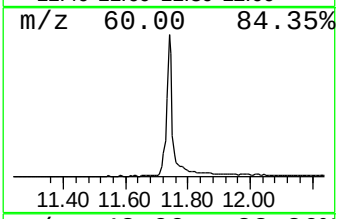
R.T.	EstConc	Area	Relative to ISTD	R.T.
11.74	16.63 ng	2743120	Phenanthrene-d10	11.19

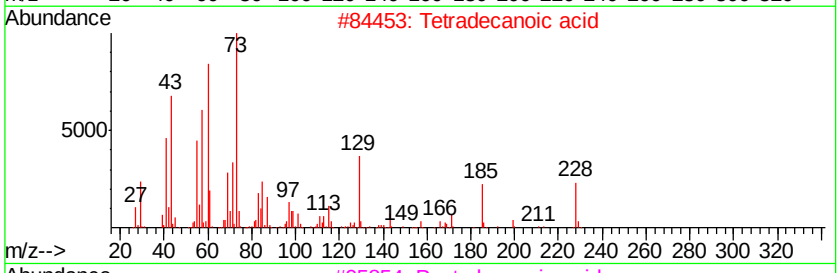
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	94
4	Octadecanoic acid	284	C18H36O2	000057-11-4	87
5	n-Decanoic acid	172	C10H20O2	000334-48-5	76

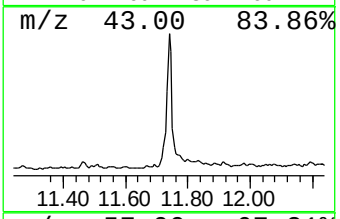


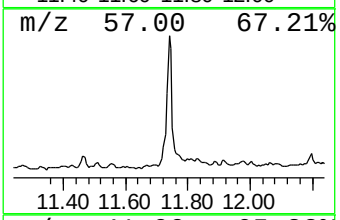


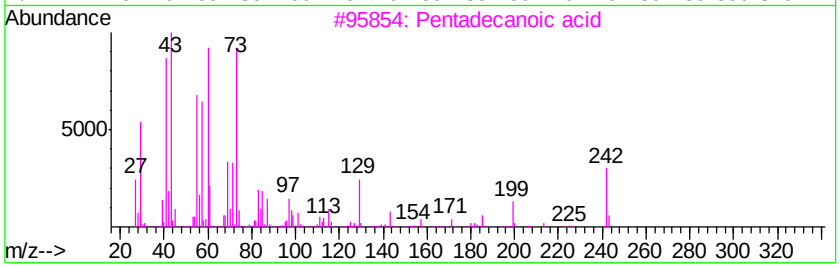


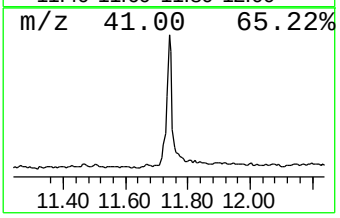












Library Search Compound Report

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Sample : K3097-23
Misc :
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Instrument : BNA_F
ClientSampleId : 0443-PR-6(PASSAIC)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

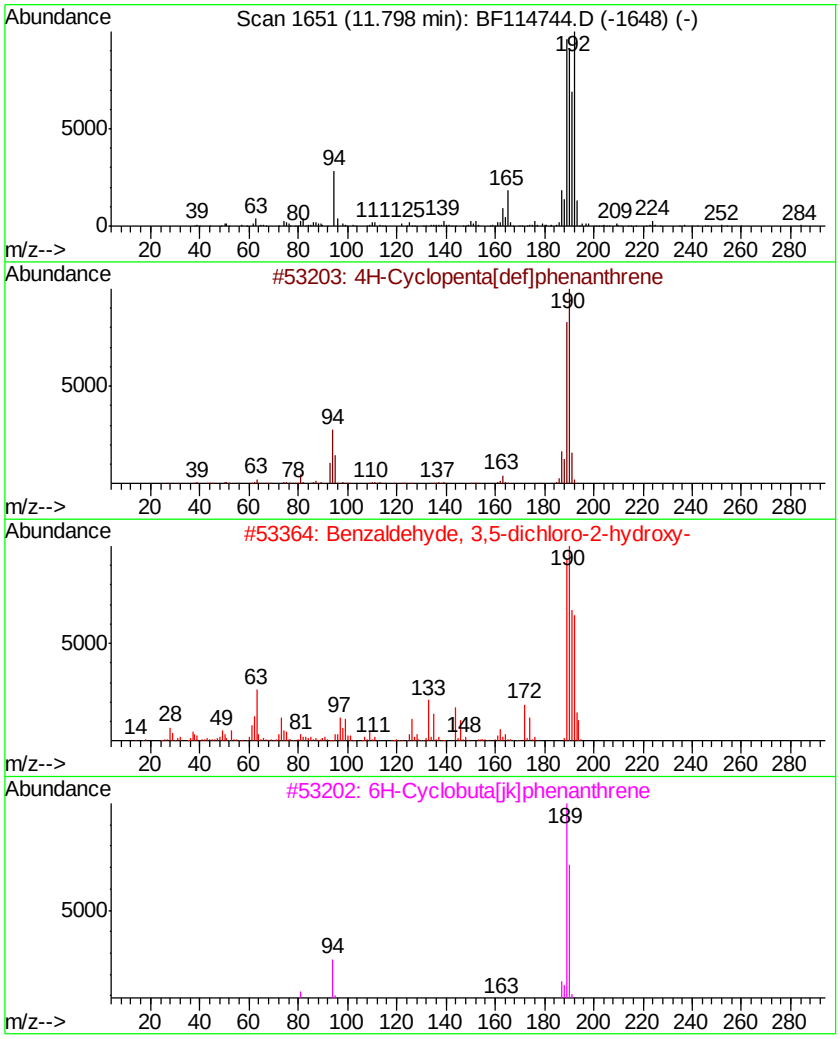
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

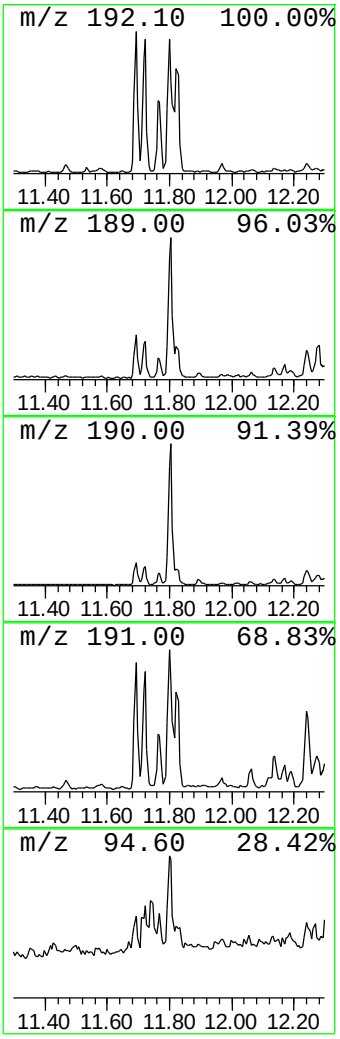
****

Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.80	2.76 ng	454636	Phenanthrene-d10	11.19

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
2	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	53
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	49
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	46
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	42





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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

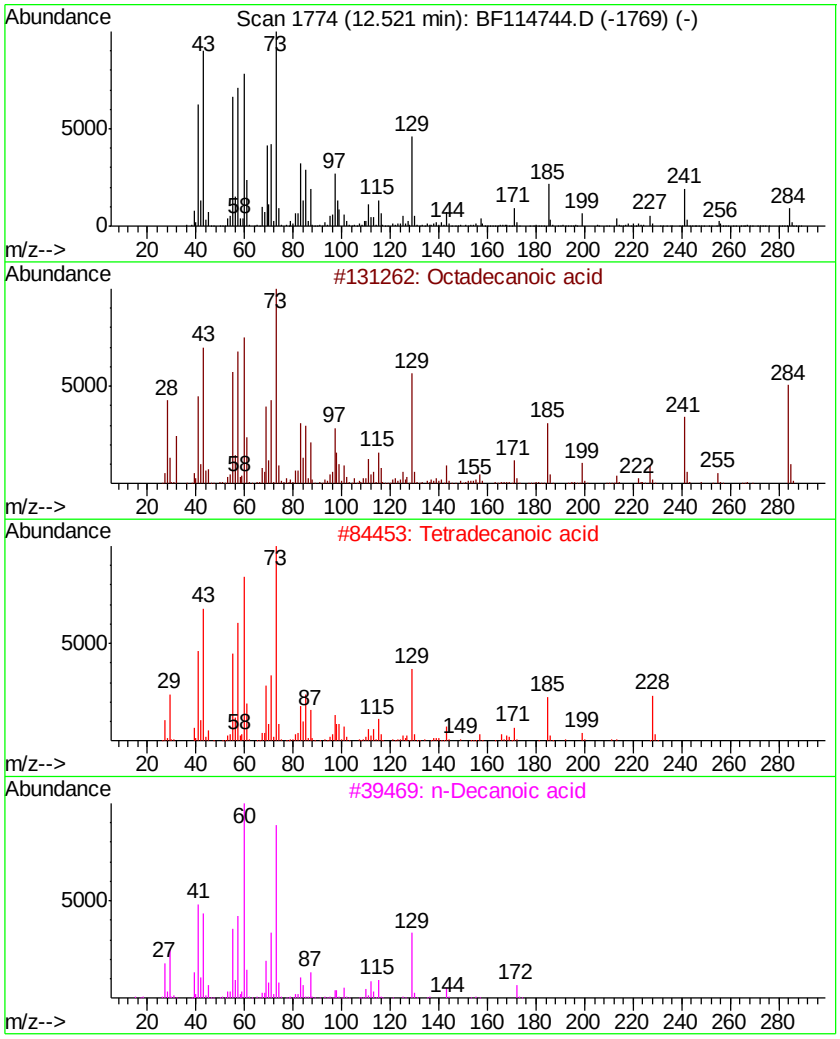
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

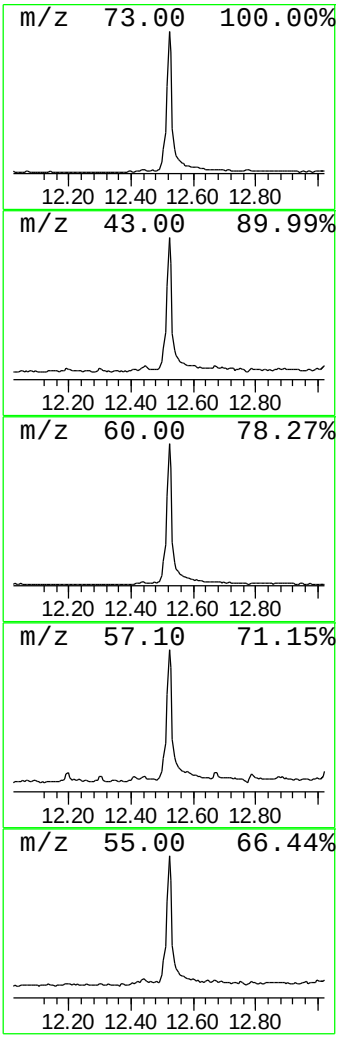
****

Peak Number 5 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.52	26.29 ng	3920630	Chrysene-d12	13.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	87
3	n-Decanoic acid	172	C10H20O2	000334-48-5	70
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	62
5	Oxalic acid, propyl tetradecyl e...	328	C19H36O4	1000309-26-7	30





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TIC Integration Parameters: LSCINT.P

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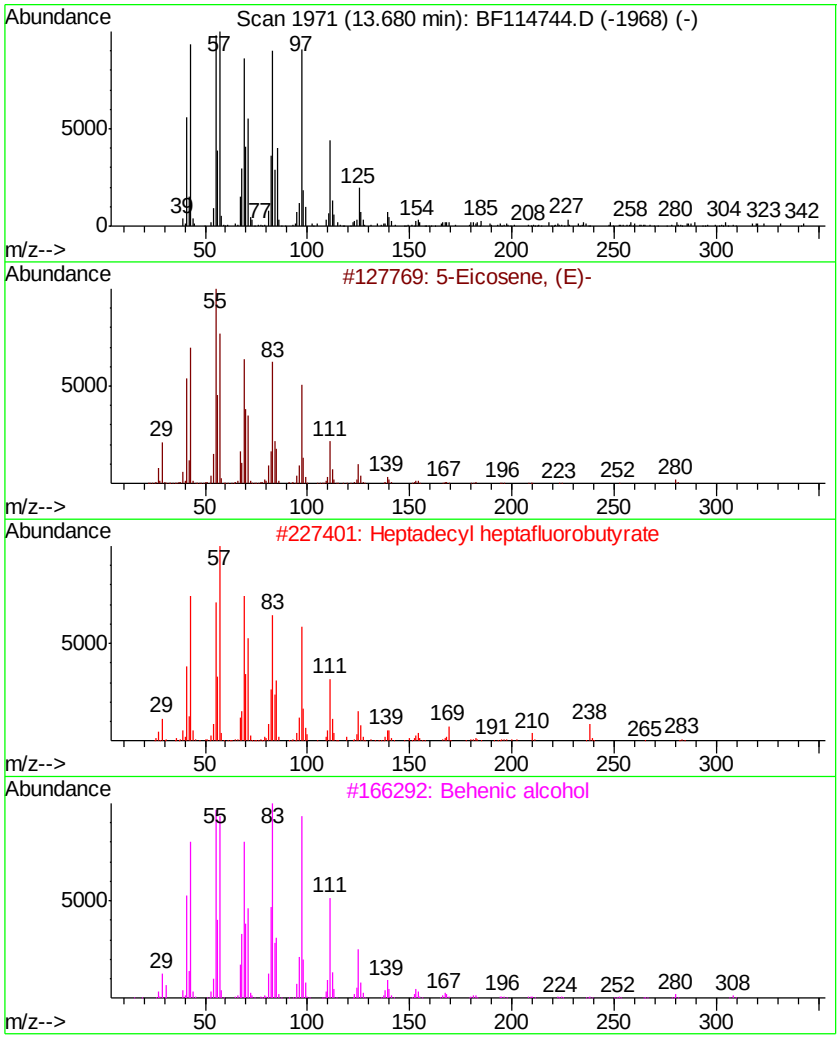
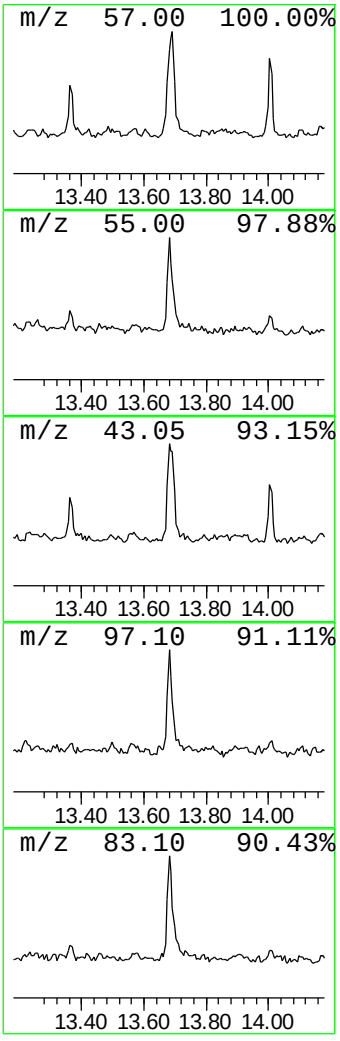
Peak Number 6 5-Eicosene, (E)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.68	3.70 ng	551121	Chrysene-d12	13.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	97
2	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93
3	Behenic alcohol	326	C22H46O	000661-19-8	93
4	3-Eicosene, (E)-	280	C20H40	074685-33-9	92
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91





Library Search Compound Report

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Data File : BF114744.D
Acq On : 1 Jun 2019 9:34
Operator : HP/JU
Sample : K3097-23
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 0443-PR-6(PASSAIC)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052919.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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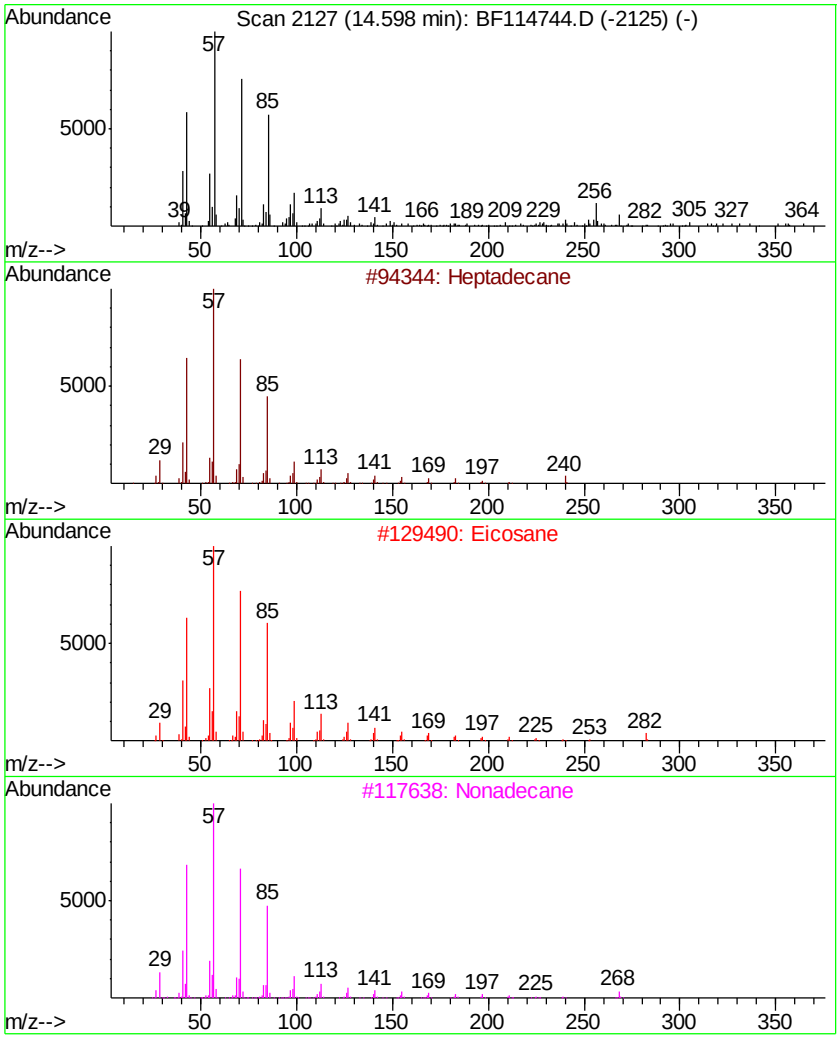
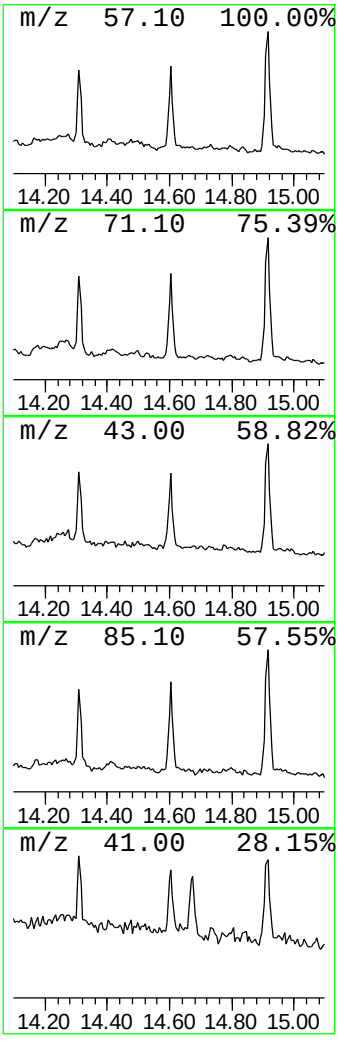
Peak Number 7 Heptadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	2.31 ng	276679	Perylene-d12	15.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	96
2	Eicosane	282	C20H42	000112-95-8	93
3	Nonadecane	268	C19H40	000629-92-5	91
4	Hexadecane	226	C16H34	000544-76-3	90
5	Heneicosane	296	C21H44	000629-94-7	87





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Sample : K3097-23
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Instrument : BNA_F
ClientSampleId : 0443-PR-6(PASSAIC)

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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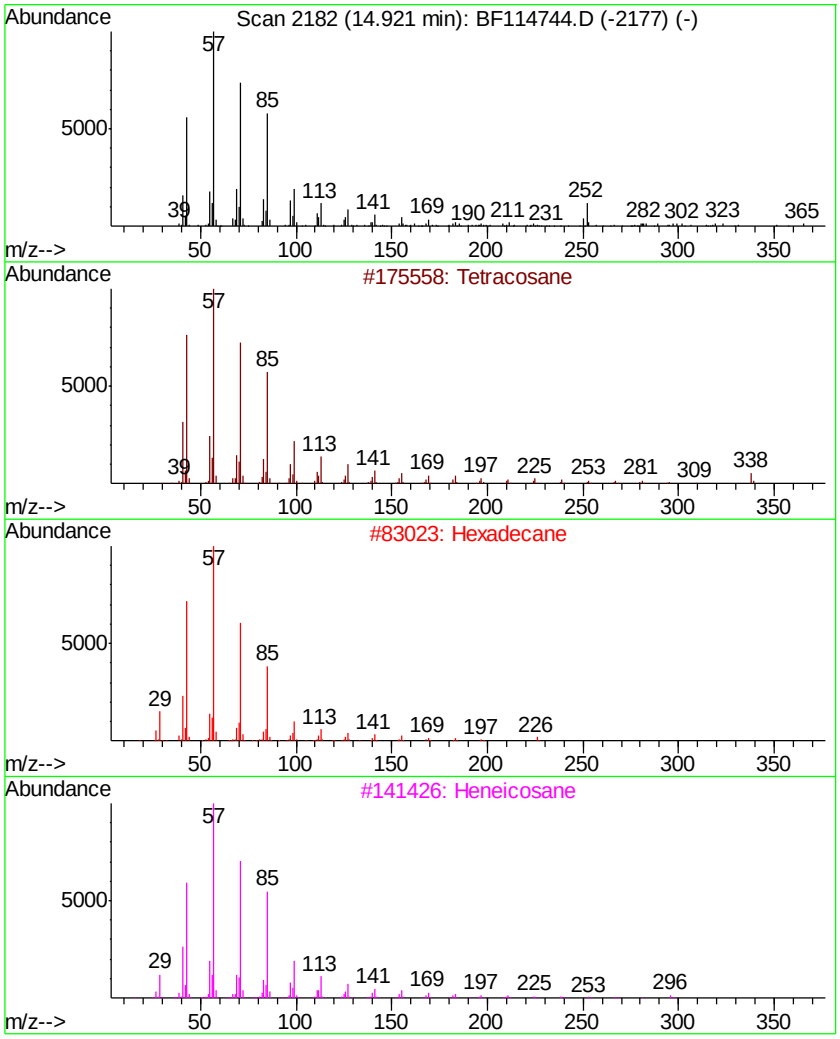
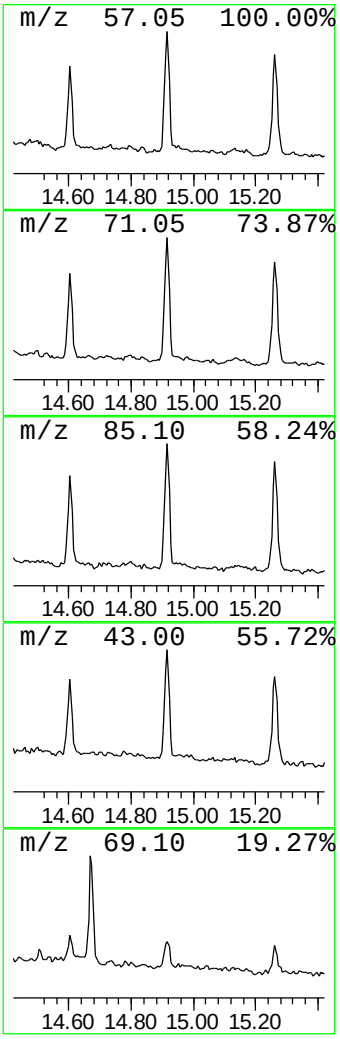
Peak Number 8 Tetracosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.92	3.70 ng	442668	Perylene-d12	15.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	97
2	Hexadecane	226	C16H34	000544-76-3	94
3	Heneicosane	296	C21H44	000629-94-7	81
4	Hexacosane	366	C26H54	000630-01-3	81
5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	76





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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 0443-PR-6(PASSAIC)

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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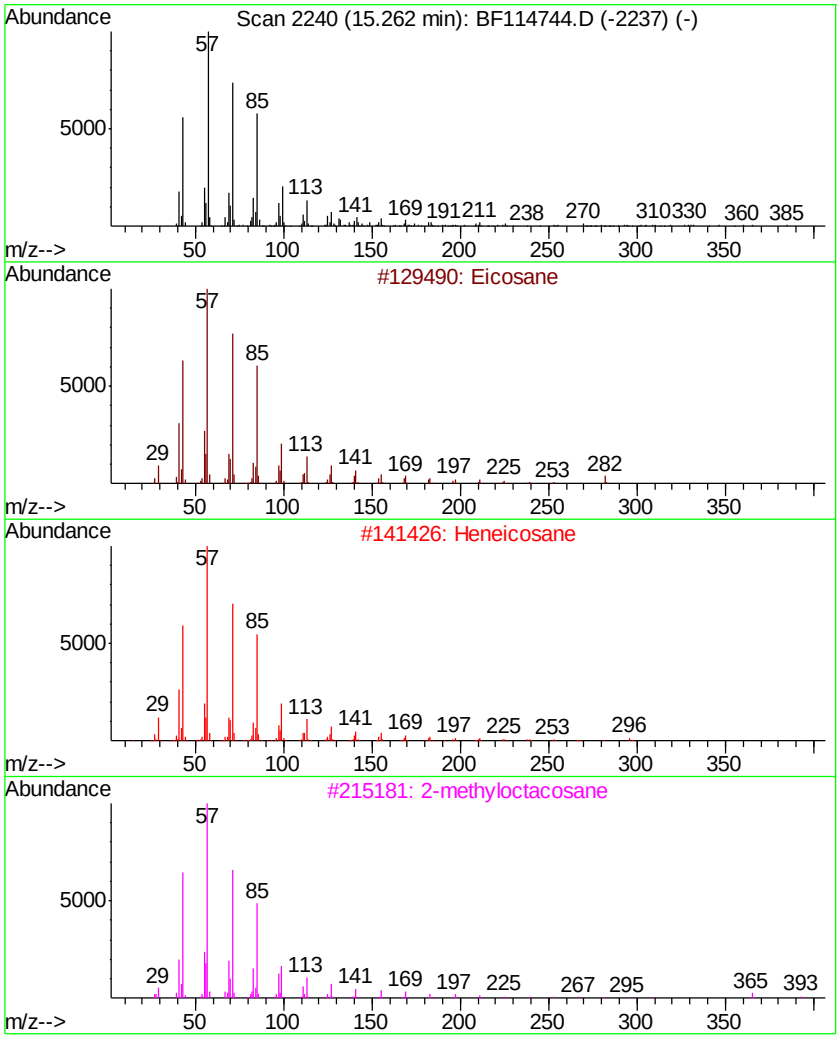
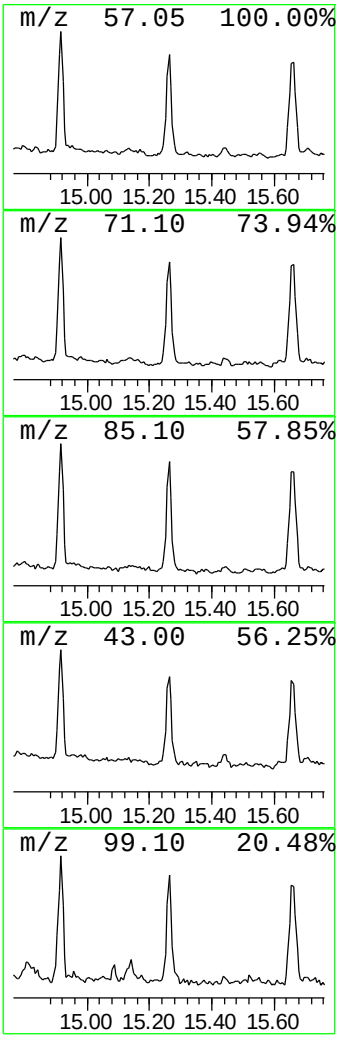
Peak Number 9 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	2.73 ng	327095	Perylene-d12	15.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	94
2	Heneicosane	296	C21H44	000629-94-7	91
3	2-methyloctacosane	408	C29H60	1000376-72-8	91
4	Hexacosane	366	C26H54	000630-01-3	87
5	Tetracosane	338	C24H50	000646-31-1	87





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Sample : K3097-23
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 0443-PR-6(PASSAIC)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052919.M
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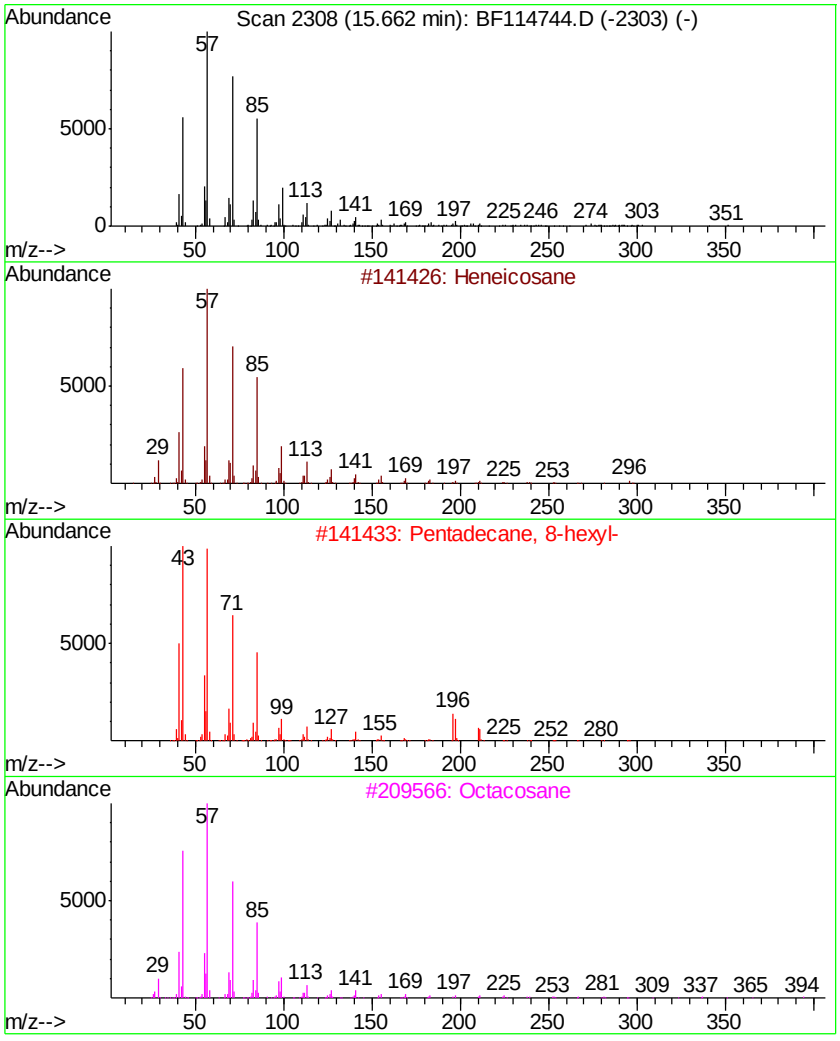
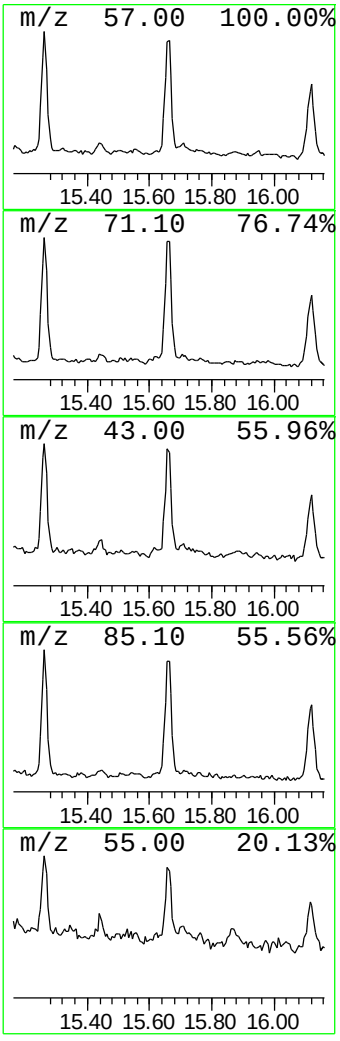
Peak Number 10 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	3.27 ng	391317	Perylene-d12	15.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
3	Octacosane	394	C28H58	000630-02-4	90
4	Eicosane, 10-methyl-	296	C21H44	054833-23-7	87
5	Heptadecane	240	C17H36	000629-78-7	87





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Acq On : 1 Jun 2019 9:34
Operator : HP/JU
Sample : K3097-23
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 0443-PR-6(PASSAIC)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052919.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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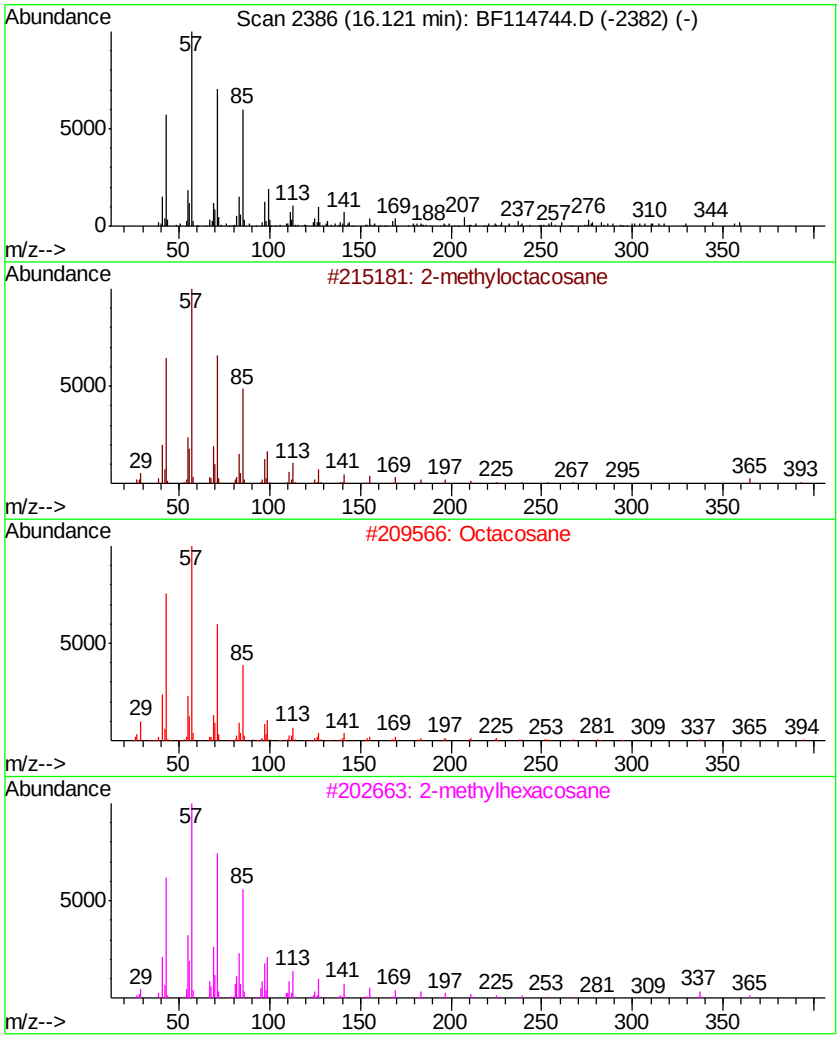
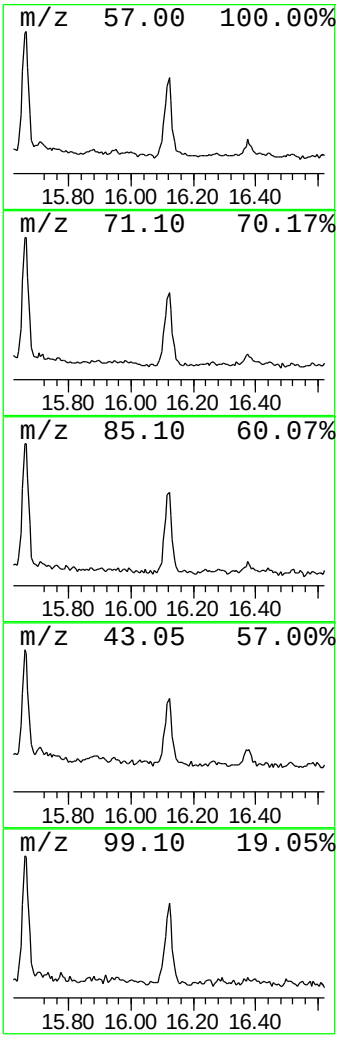
Peak Number 11 2-methyloctacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.12	2.66 ng	318689	Perylene-d12	15.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-methyloctacosane	408	C29H60	1000376-72-8	86
2	Octacosane	394	C28H58	000630-02-4	86
3	2-methylhexacosane	380	C27H56	1000376-72-7	86
4	Octadecane	254	C18H38	000593-45-3	83
5	Docosane	310	C22H46	000629-97-0	83





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Operator : HP/JU
Sample : K3097-23
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 0443-PR-6(PASSAIC)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052919.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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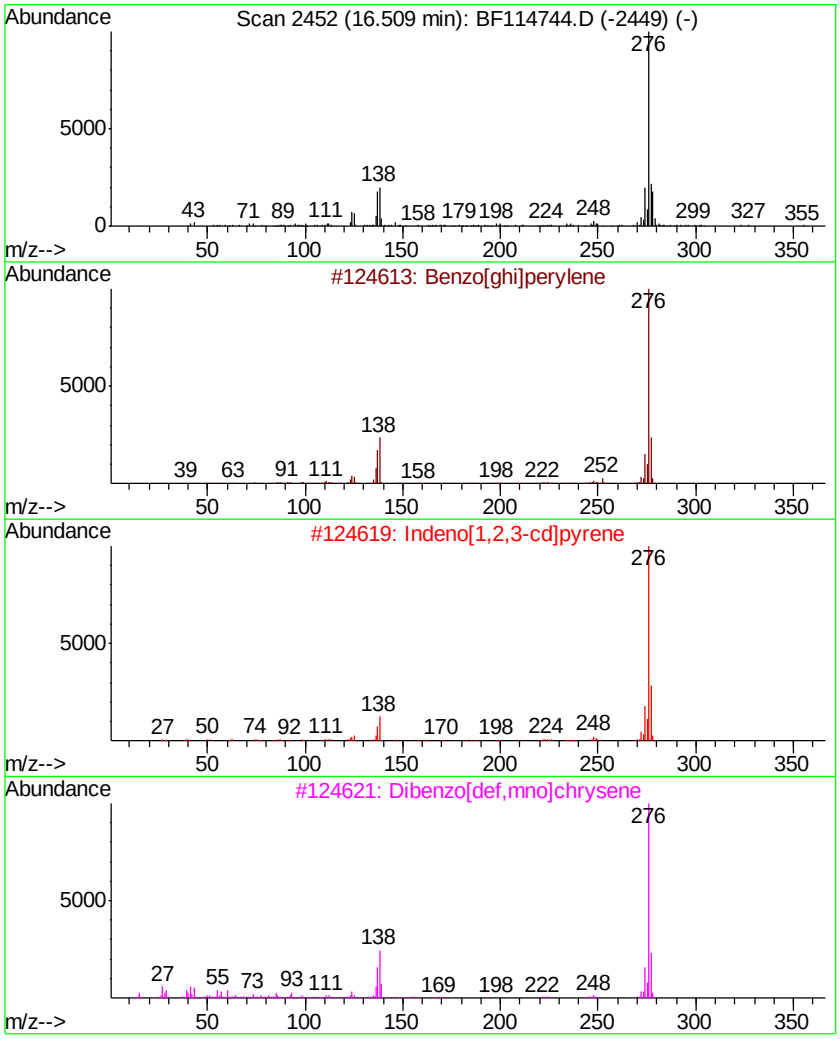
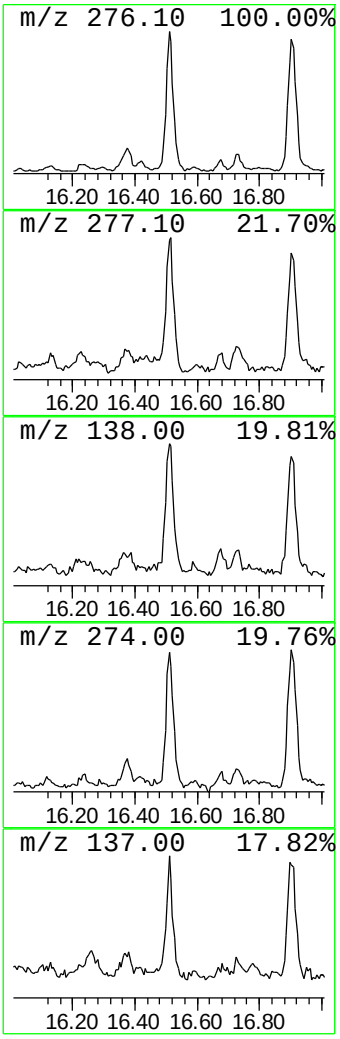
Peak Number 12 Dibenzo[def,mno]chrysene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.51	2.42 ng	290339	Perylene-d12	15.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[ghi]perylene	276	C22H12	000191-24-2	95
2	Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	93
3	Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	90
4	6H-Pyrido[4,3-b]carbazole, 9-met...	276	C18H16N2O	010371-86-5	47
5	1-Naphthalenecarboxylic acid, 2-...	276	C18H12O3	038119-11-8	43





Tentatively Identified Compound (LSC) summary

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Data File : BF114744.D
Acq On : 1 Jun 2019 9:34
Operator : HP/JU
Sample : K3097-23
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 0443-PR-6(PASSAIC)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052919.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
unknown6.47	6.47	85.4	ng	9285040	1	6.69	2173710	20.0
n-Hexadecanoic acid	11.74	16.6	ng	2743120	4	11.19	3299580	20.0
4H-Cyclopenta[def...	11.80	2.8	ng	454636	4	11.19	3299580	20.0
Octadecanoic acid	12.52	26.3	ng	3920630	5	13.81	2982500	20.0
5-Eicosene, (E)-	13.68	3.7	ng	551121	5	13.81	2982500	20.0
Heptadecane	14.60	2.3	ng	276679	6	15.20	2395570	20.0
Tetracosane	14.92	3.7	ng	442668	6	15.20	2395570	20.0
Eicosane	15.26	2.7	ng	327095	6	15.20	2395570	20.0
Heneicosane	15.66	3.3	ng	391317	6	15.20	2395570	20.0
2-methyloctacosane	16.12	2.7	ng	318689	6	15.20	2395570	20.0
Dibenzo[def,mno]c...	16.51	2.4	ng	290339	6	15.20	2395570	20.0