LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041117\
Data File : BF094373.D
Acq On : 12 Apr 2017 1:50
Operator : SJ/MA
Sample : I2610-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SOURCE3-SE

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:\HPCHEM1\BNA_F\METHODS\8270-BF032217.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	3.945	347	350	357	rBV	37565	38129	3.36%	0.259%
2	4.375	419	423	435	rBV	446788	465122	41.01%	3.161%
3	5.457	603	607	617	rVB	584968	586466	51.71%	3.985%
4	5.581	624	628	635	rVB	589273	539848	47.60%	3.668%
5	5.828	666	670	674	rBV	963917	848216	74.80%	5.764%
6	6.004	696	700	704	rBV	544202	473139	41.72%	3.215%
7	6.469	775	779	784	rBV	357455	349635	30.83%	2.376%
8	7.075	878	882	887	rVB7	9372	16794	1.48%	0.114%
9	7.333	920	926	929	rBV	1178972	1134041	100.00%	7.706%
10	7.357	929	930	934	rVB	86633	46455	4.10%	0.316%
11	8.098	1050	1056	1058	rBV3	10032	14426	1.27%	0.098%
12	8.198	1069	1073	1080	rBV	90814	88581	7.81%	0.602%
13	8.316	1088	1093	1096	rBV	73899	69300	6.11%	0.471%
14	8.651	1146	1150	1154	rVB	698092	621405	54.80%	4.222%
15	8.769	1168	1170	1172	rBV	14117	11833	1.04%	0.080%
16	8.792	1172	1174	1180	rVB3	12377	12664	1.12%	0.086%
17	8.886	1188	1190	1195	rVB	18356	18111	1.60%	0.123%
18	8.969	1200	1204	1209	rBV2	30159	49677	4.38%	0.338%
19	9.063	1216	1220	1222	rBV	49163	52040	4.59%	0.354%
20	9.092	1222	1225	1227	rVV	32226	25450	2.24%	0.173%
21	9.151	1231	1235	1239	rVV2	61352	55865	4.93%	0.380%
22	9.204	1239	1244	1246	rBV2	17206	22497	1.98%	0.153%
23	9.298	1254	1260	1265	rVB2	66333	81951	7.23%	0.557%
24	9.469	1282	1289	1294	rBV2	1001268	1105327	97.47%	7.511%
25	9.510	1294	1296	1298	rBV	22575	16185	1.43%	0.110%
26	9.616	1310	1314	1319	rBV6	12092	18130	1.60%	0.123%
27	9.727	1329	1333	1338	rVB4	18406	31062	2.74%	0.211%
28	9.780	1338	1342	1347	rBV	19885	26937	2.38%	0.183%
29	9.874	1354	1358	1362	rBV3	14888	25161	2.22%	0.171%
30	9.980	1372	1376	1381	rBV6	14069	25613	2.26%	0.174%
31	10.063	1387	1390	1393	rVB	22071	18919	1.67%	0.129%
32	10.110	1396	1398	1401	rVB	16180	13987	1.23%	0.095%
33	10.145	1401	1404	1409	rBV2	56105	54458	4.80%	0.370%
34	10.339	1434	1437	1440	rBV3	12712	14695	1.30%	0.100%

LSC Area Percent Report

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Acq On : 12 Apr 2017 1:50
Operator : SJ/MA
Sample : I2610-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SOURCE3-SE

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:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	10.445	1451	1455	1462	rVB	250150	241288	21.28%	1.640%
36	10.598	1477	1481	1486	rVB	122231	126503	11.16%	0.860%
37	10.651	1486	1490	1491	rBV	29709	29973	2.64%	0.204%
38	10.821	1515	1519	1522	rVB4	17408	25915	2.29%	0.176%
39	10.857	1522	1525	1528	rVB2	16064	15405	1.36%	0.105%
40	10.898	1529	1532	1535	rBV5	17469	17887	1.58%	0.122%
41	10.927	1535	1537	1540	rVB3	20316	16566	1.46%	0.113%
42	10.969	1540	1544	1545	rBV4	9675	12223	1.08%	0.083%
43	11.057	1554	1559	1562	rBV3	11905	16402	1.45%	0.111%
44	11.169	1575	1578	1581	rBV	32709	37157	3.28%	0.252%
45	11.204	1581	1584	1587	rVB2	22718	21357	1.88%	0.145%
46	11.245	1587	1591	1595	rBV7	10909	17009	1.50%	0.116%
47	11.298	1595	1600	1602	rBV	990595	958713	84.54%	6.514%
48	11.321	1602	1604	1608	rVV	264517	243882	21.51%	1.657%
49	11.386	1612	1615	1619	rVB	46911	46452	4.10%	0.316%
50	11.451	1619	1626	1628	rBV6	11575	21283	1.88%	0.145%
51	11.704	1666	1669	1672	rBV	28753	27993	2.47%	0.190%
52	11.733	1672	1674	1677	rVB2	14530	12835	1.13%	0.087%
53	11.810	1685	1687	1691	rVB	19103	19057	1.68%	0.129%
54	11.857	1691	1695	1699	rBV5	29110	39675	3.50%	0.270%
55	11.921	1702	1706	1709	rBV	73557	78704	6.94%	0.535%
56	11.957	1709	1712	1716	rVB	81497	85708	7.56%	0.582%
57	12.027	1718	1724	1725	rBV2	51723	70974	6.26%	0.482%
58	12.051	1725	1728	1731	rVV3	110798	132548	11.69%	0.901%
59	12.080	1731	1733	1736	rVB	50103	48261	4.26%	0.328%
60	12.239	1757	1760	1762	rBV3	12998	13616	1.20%	0.093%
61	12.292	1766	1769	1771	rBV2	31910	32295	2.85%	0.219%
62	12.486	1798	1802	1804	rBV3	64633	98965	8.73%	0.672%
63	12.604	1818	1822	1824	rBV2	60628	76733	6.77%	0.521%
64	12.639	1824	1828	1831	rVV3	179483	210638	18.57%	1.431%
65	12.698	1835	1838	1839	rVV	70257	72704	6.41%	0.494%
66	12.721	1839	1842	1844	rVV3	121196	146011	12.88%	0.992%
67	12.757	1844	1848	1851	rVV3	185500	265020	23.37%	1.801%
68	12.786	1851	1853	1857	rVB	192382	189431	16.70%	1.287%
69	12.868	1864	1867	1870	rBV4	21104	28840	2.54%	0.196%
70	12.910	1870	1874	1875	rBV2	87894	115392	10.18%	0.784%
71	12.927	1875	1877	1881	rVB2	127499	136954	12.08%	0.931%

LSC Area Percent Report

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Acq On : 12 Apr 2017 1:50
Operator : SJ/MA
Sample : I2610-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SOURCE3-SE

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:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	13.063	1896	1900	1904	rBV	281766	290534	25.62%	1.974%
73	13.274	1931	1936	1939	rBV	437432	486245	42.88%	3.304%
74	13.321	1939	1944	1947	rVV	181937	230245	20.30%	1.565%
75	13.604	1990	1992	2001	rVB3	51583	66657	5.88%	0.453%
76	13.715	2009	2011	2014	rBV	40508	36262	3.20%	0.246%
77	13.757	2015	2018	2023	rBV4	78555	144799	12.77%	0.984%
78	13.910	2039	2044	2050	rBV	189368	269629	23.78%	1.832%
79	14.139	2079	2083	2089	rVB2	177250	244491	21.56%	1.661%
80	14.551	2148	2153	2157	rBV	816509	1035992	91.35%	7.040%
81	16.180	2427	2430	2439	rVB	570732	661981	58.37%	4.498%
82	16.880	2546	2549	2557	rVB4	344678	527310	46.50%	3.583%

Sum of corrected areas: 14716633

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Operator : SJ/MA
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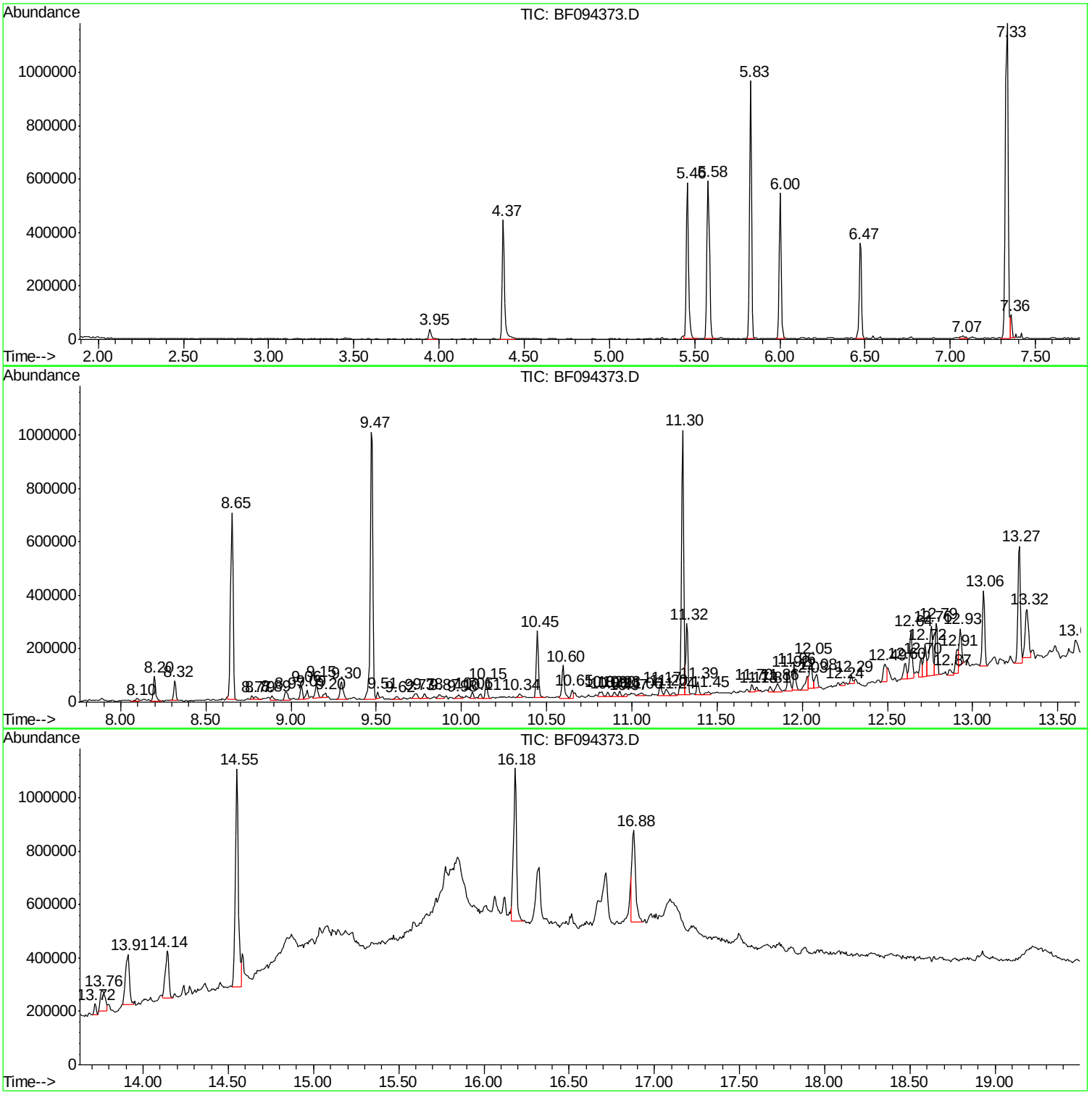
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SOURCE3-SE

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\HPCHEM1\BNA F\DATA\BF041117\
Data File : BF094373.D
Acq On : 12 Apr 2017 1:50
Operator : SJ/MA
Sample : I2610-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SOURCE3-SE

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

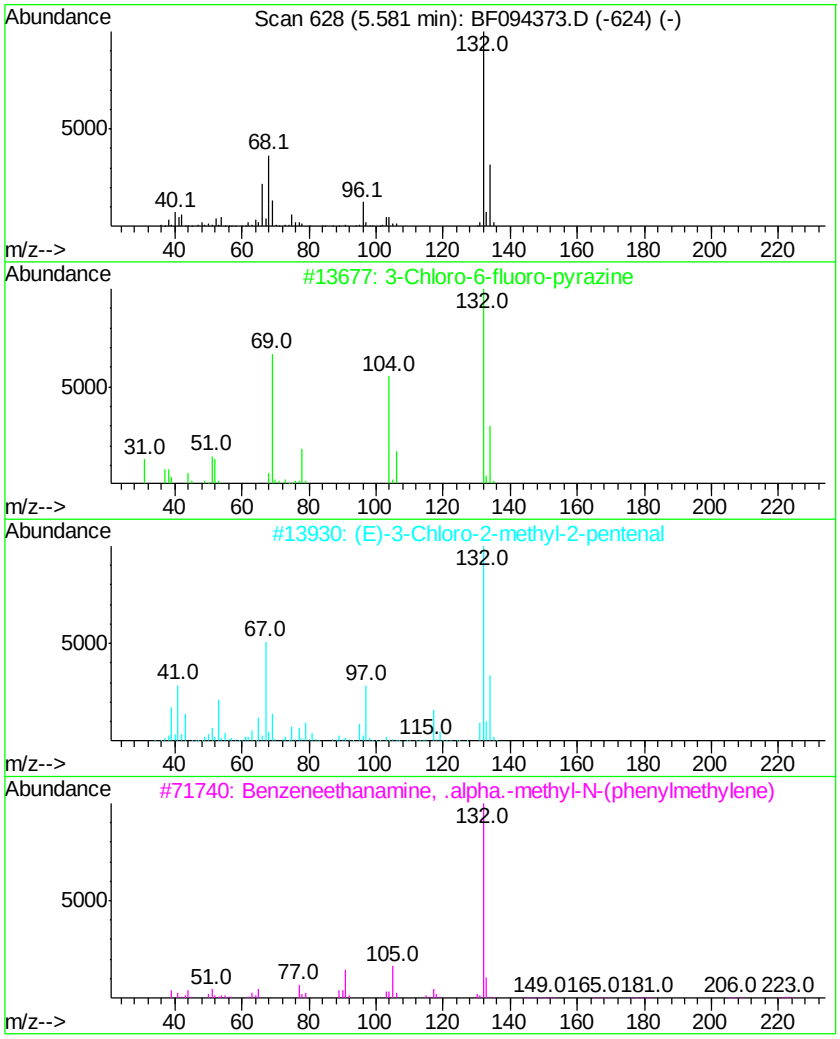
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

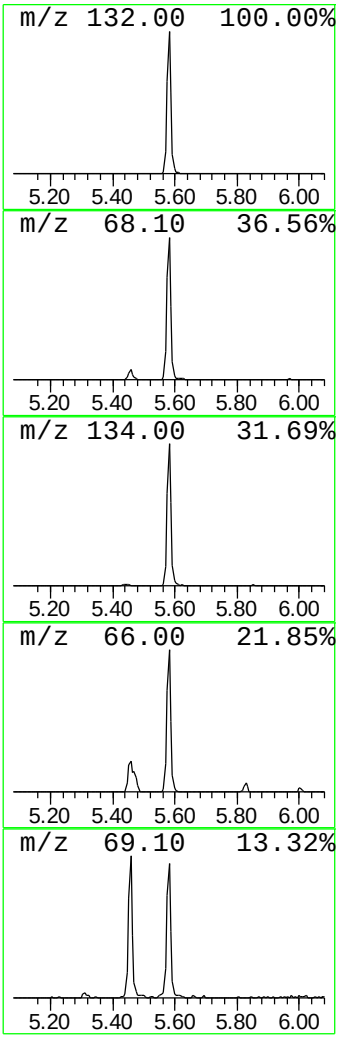
****

Peak Number 1 unknown5.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.58	12.73 ng	539848	1,4-Dichlorobenzene-d4	5.83

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	Benzeneethanamine, .alpha.-methy...	223	C16H17N	002980-02-1	12
4	5-Aminoindole	132	C8H8N2	005192-03-0	10
5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9





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Data File : BF094373.D
Acq On : 12 Apr 2017 1:50
Operator : SJ/MA
Sample : I2610-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SOURCE3-SE

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

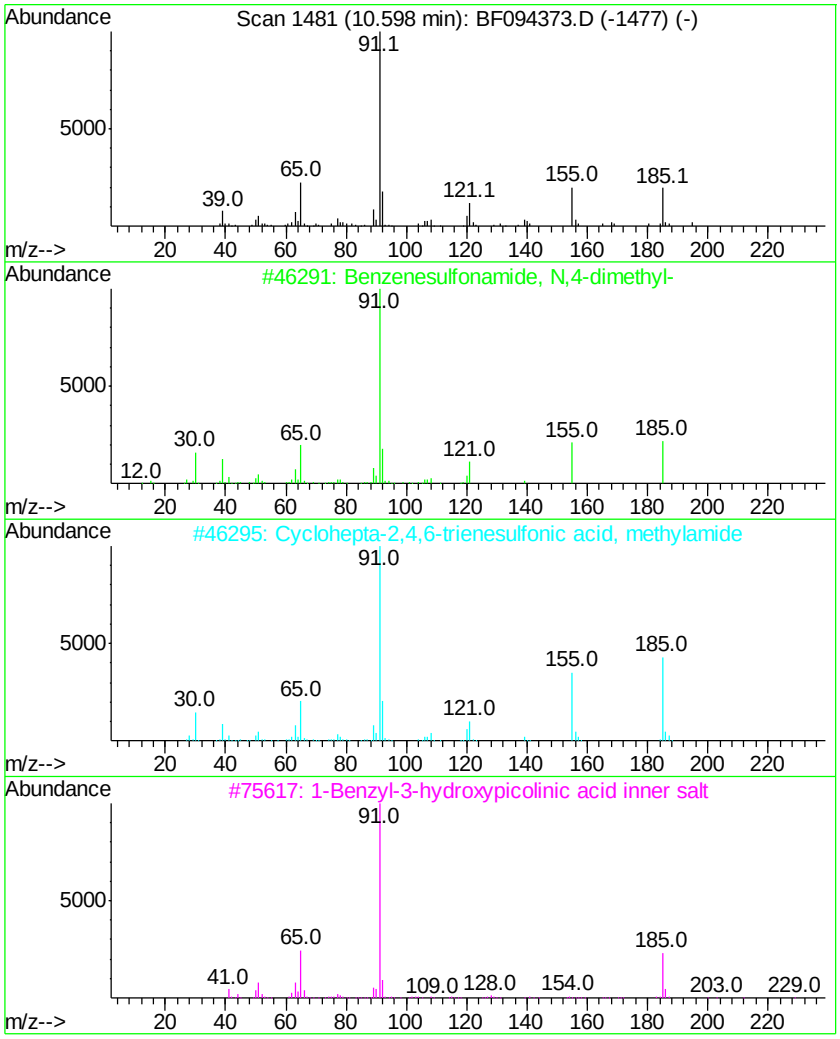
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

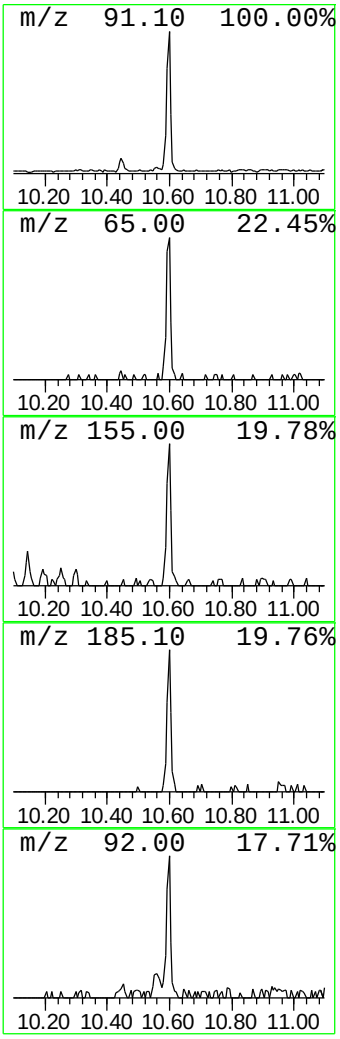
****

Peak Number 3 Benzenesulfonamide, N,4-dim... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.60	2.64 ng	126503	Phenanthrene-d10	11.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N,4-dimethyl-	185	C8H11NO2S	000640-61-9	94
2	Cyclohepta-2,4,6-trienesulfonic ...	185	C8H11NO2S	1000187-28-7	49
3	1-Benzyl-3-hydroxypicolinic acid...	229	C13H11NO3	1000212-09-1	47
4	p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	47
5	(O-p-Tosylisonitroso)malononitrile	249	C10H7N3O3S	020893-01-0	47





Data Path : Z:\HPCHEM1\BNA F\DATA\BF041117\
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Sample : I2610-01 5X
Misc :
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Instrument : BNA_F
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
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TIC Library : C:\DATABASE\NIST02.L
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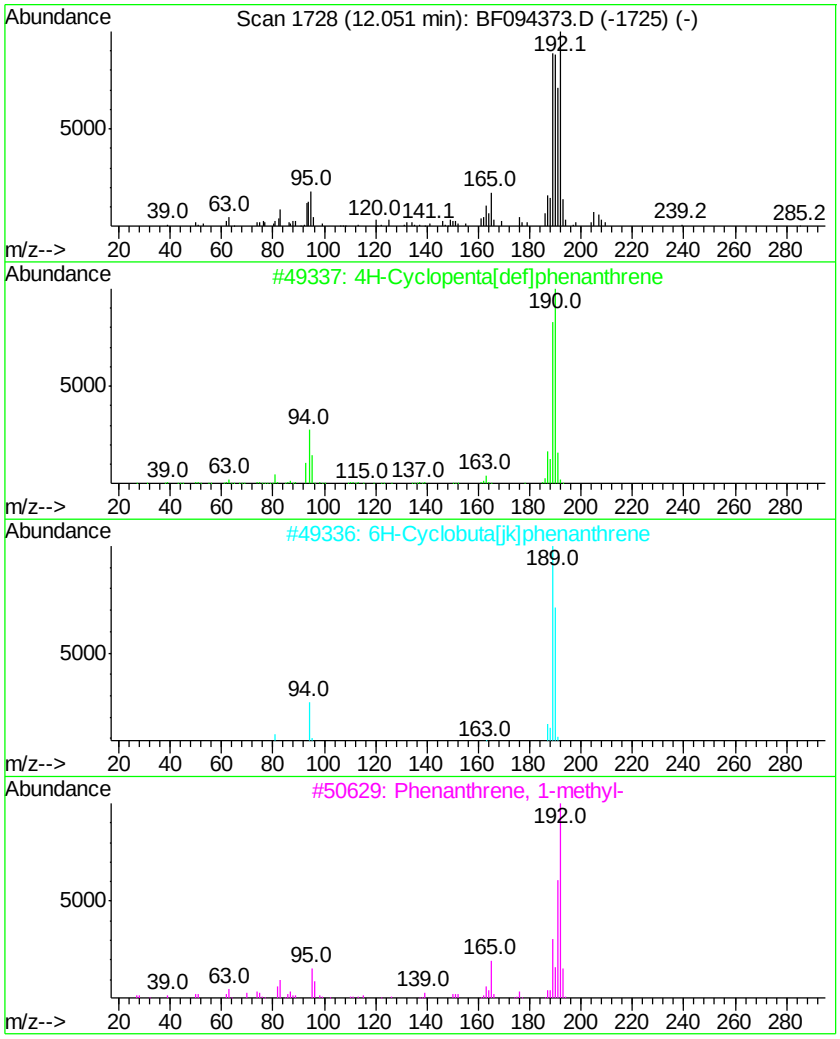
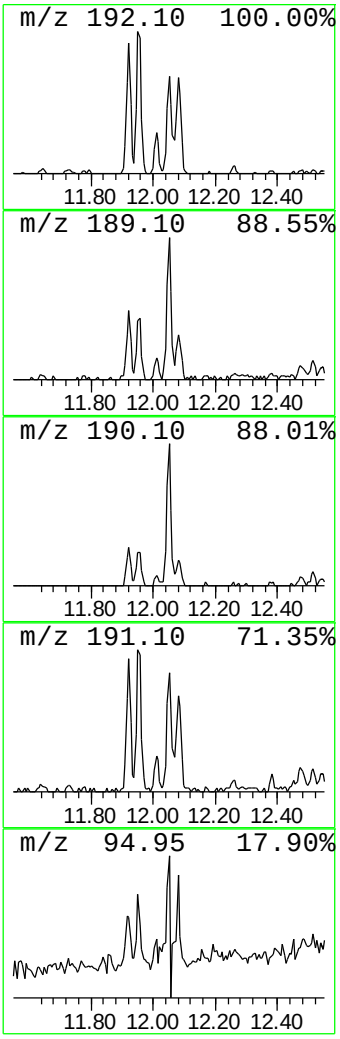
****

Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.05	2.77 ng	132548	Phenanthrene-d10	11.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	46
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	41
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	41
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	38





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TIC Library : C:\DATABASE\NIST02.L
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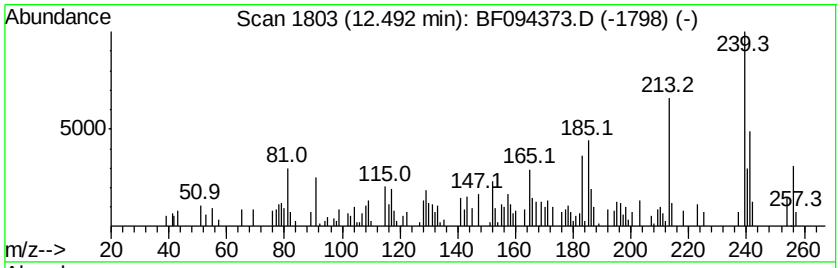
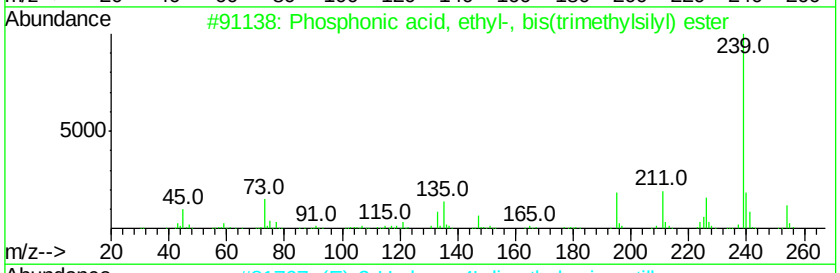
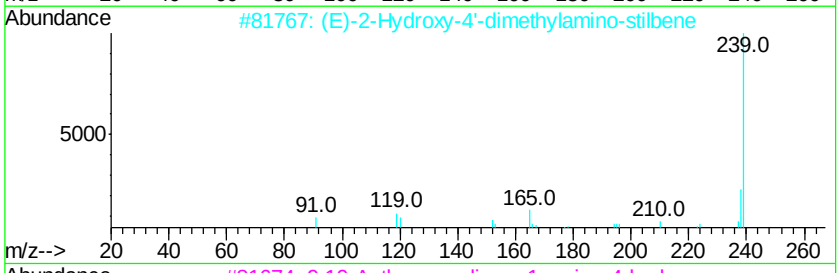
Peak Number 5 Phosphonic acid, ethyl-, bi... Concentration Rank 13

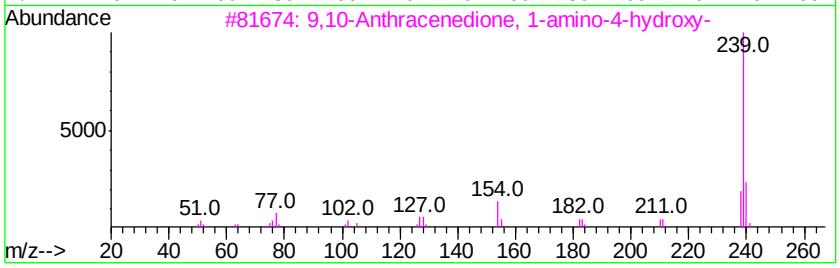
R.T.	EstConc	Area	Relative to ISTD	R.T.
12.49	2.06 ng	98965	Phenanthrene-d10	11.30

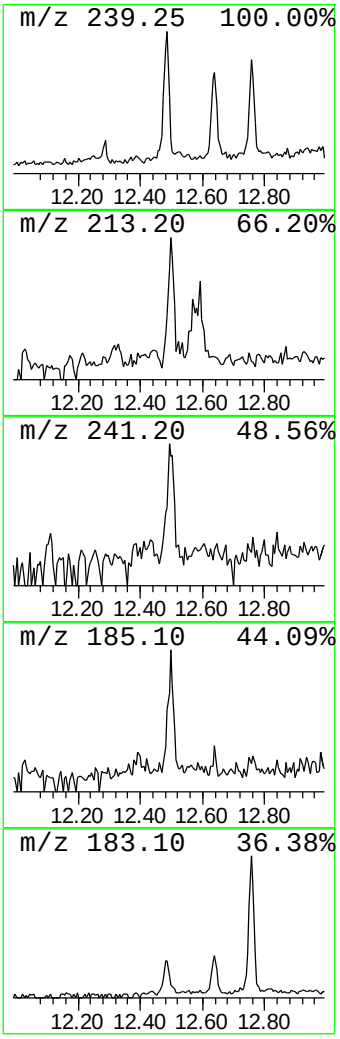
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phosphonic acid, ethyl-, bis(tri...	254	C8H23O3PSi2	001641-57-2	64
2	(E)-2-Hydroxy-4'-dimethylamino-s...	239	C16H17NO	110983-42-1	53
3	9,10-Anthracenedione, 1-amino-4-...	239	C14H9NO3	000116-85-8	53
4	p-Dimethylaminobenzylidene p-ani...	254	C16H18N2O	001749-04-8	49
5	3,6-Bis(N-methylamino)-9-methylc...	239	C15H17N3	098786-00-6	49











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Data File : BF094373.D
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Sample : I2610-01 5X
Misc :
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Instrument : BNA_F
ClientSampleId : SOURCE3-SE

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

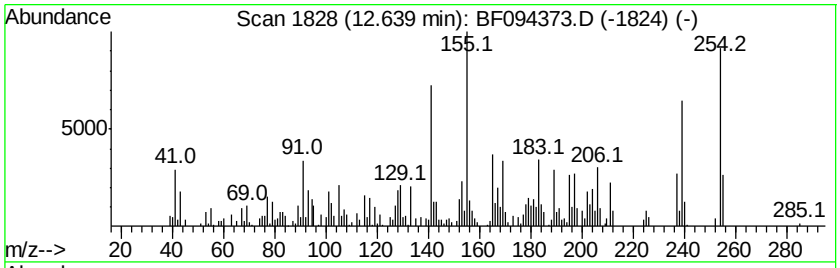
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

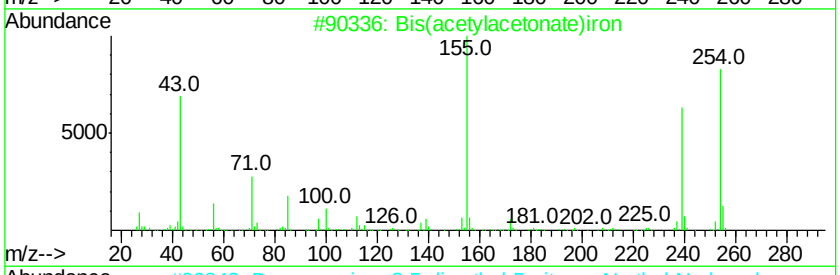
****

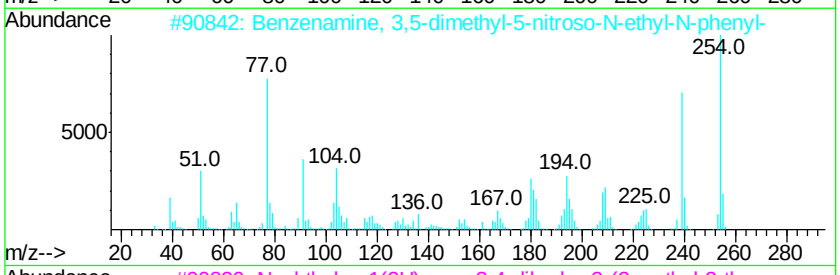
Peak Number 6 Bis(acetylacetonate)iron Concentration Rank 6

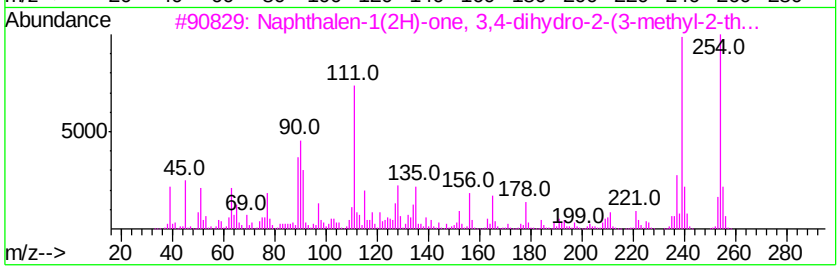
R.T.	EstConc	Area	Relative to ISTD	R.T.
12.64	4.39 ng	210638	Phenanthrene-d10	11.30

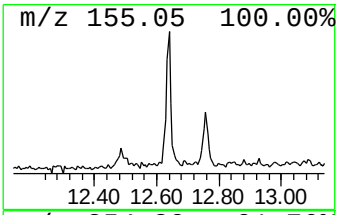
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bis(acetylacetonate)iron	254	C10H14FeO4	014024-17-0	52
2	Benzenamine, 3,5-dimethyl-5-nitr...	254	C16H18N2O	295780-40-4	42
3	Naphthalen-1(2H)-one, 3,4-dihydr...	254	C16H14OS	351066-46-1	27
4	Tetrazol-5-amine, 1-(2-methyl-1-...	239	C13H13N5	311799-88-9	25
5	(6-Isopropyl-3,4-bis(methylamino...	254	C15H18N4	014203-76-0	25

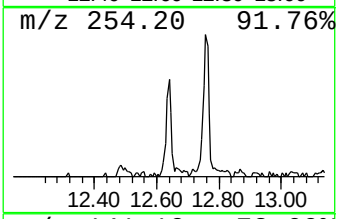


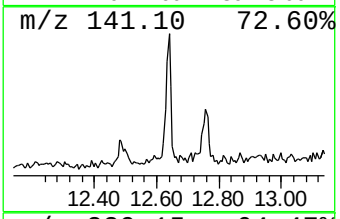


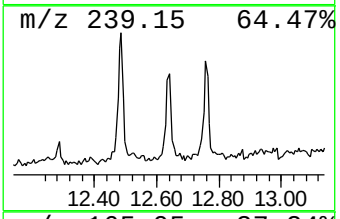


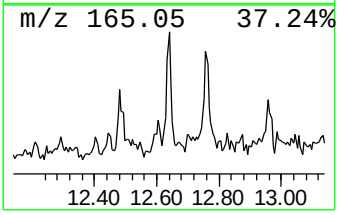












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Data File : BF094373.D
Acq On : 12 Apr 2017 1:50
Operator : SJ/MA
Sample : I2610-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SOURCE3-SE

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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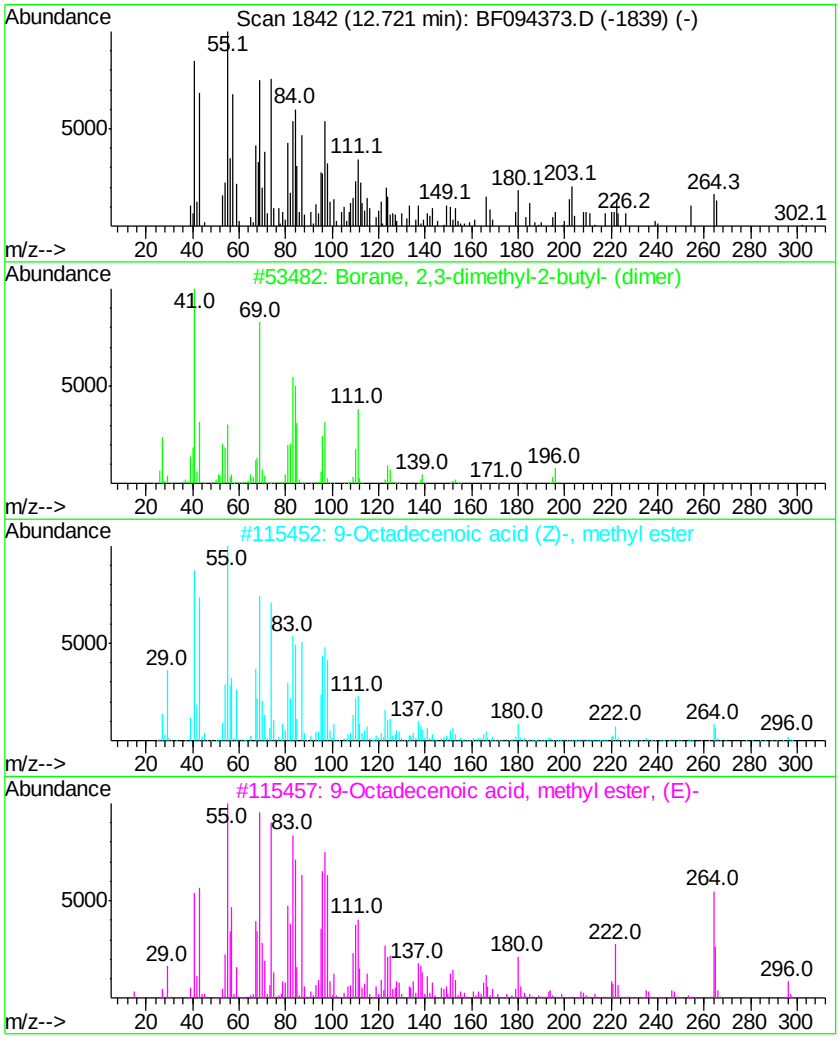
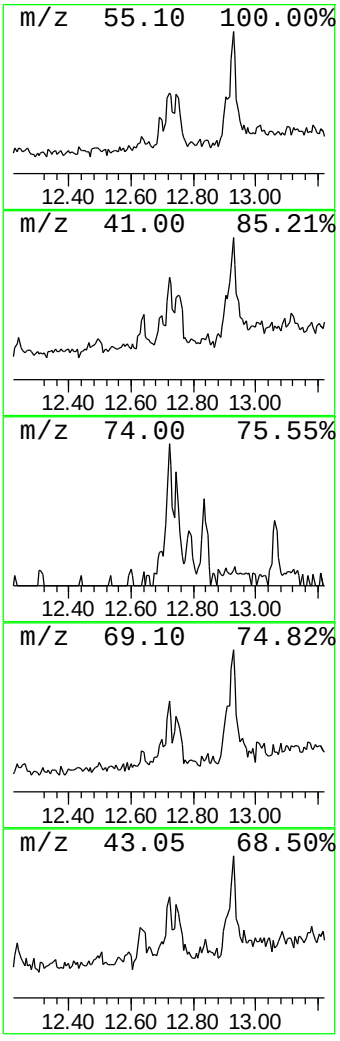
Peak Number 7 Borane, 2,3-dimethyl-2-buty... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.72	3.05 ng	146011	Phenanthrene-d10	11.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Borane, 2,3-dimethyl-2-butyl- (d...	196	C12H30B2	1000159-64-3	93
2	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	87
3	9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	87
4	9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	58
5	10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	58





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Sample : I2610-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SOURCE3-SE

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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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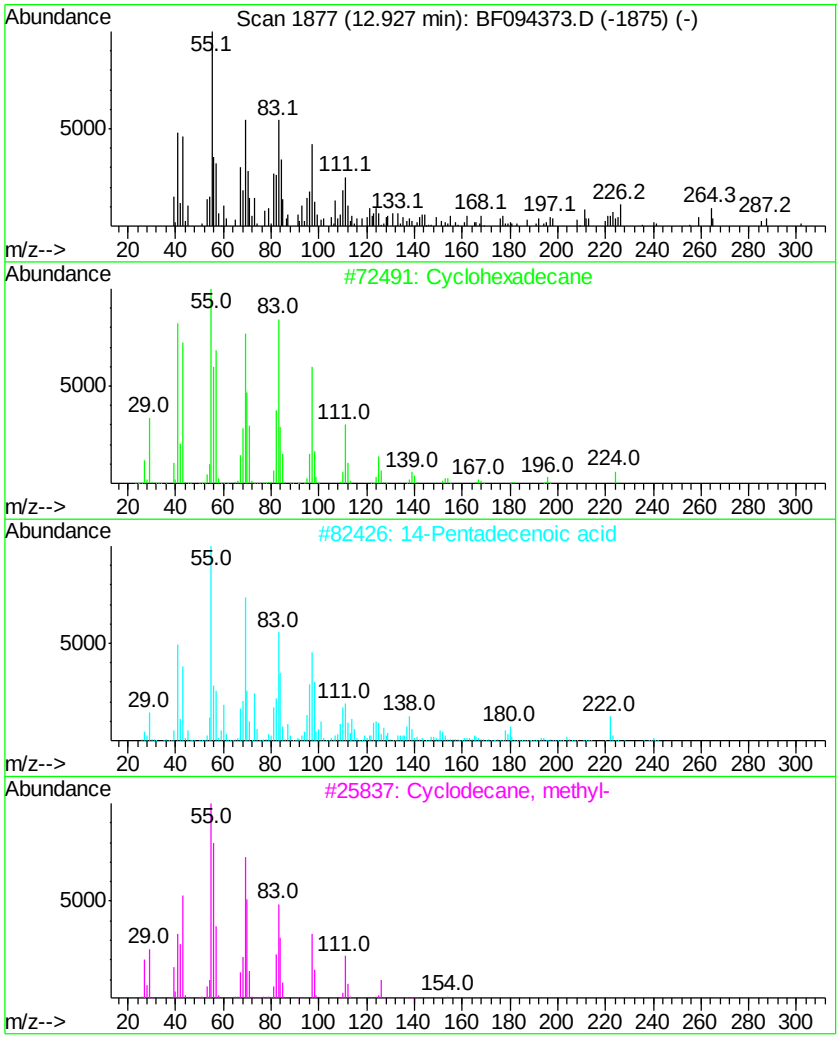
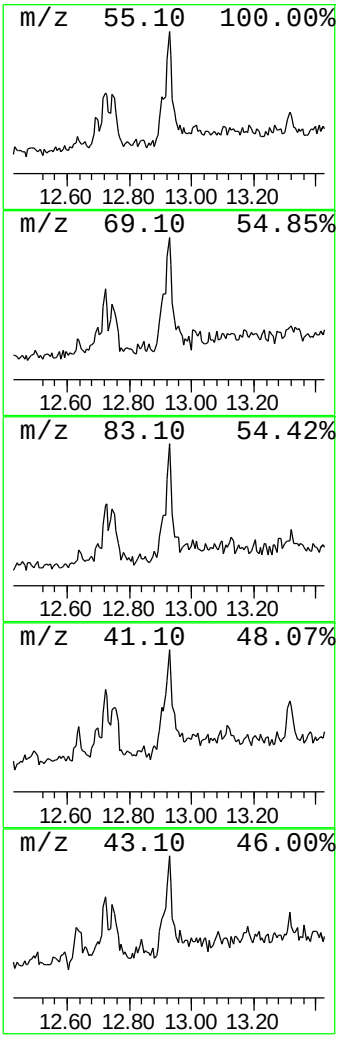
Peak Number 9 Cyclohexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.93	2.64 ng	136954	Chrysene-d12	14.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	96
2	14-Pentadecenoic acid	240	C15H28O2	017351-34-7	64
3	Cyclodecane, methyl-	154	C11H22	013151-43-4	64
4	Z-11-Pentadecenol	226	C15H30O	1000130-77-0	58
5	Cyclopentadecane	210	C15H30	000295-48-7	58





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Sample : I2610-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SOURCE3-SE

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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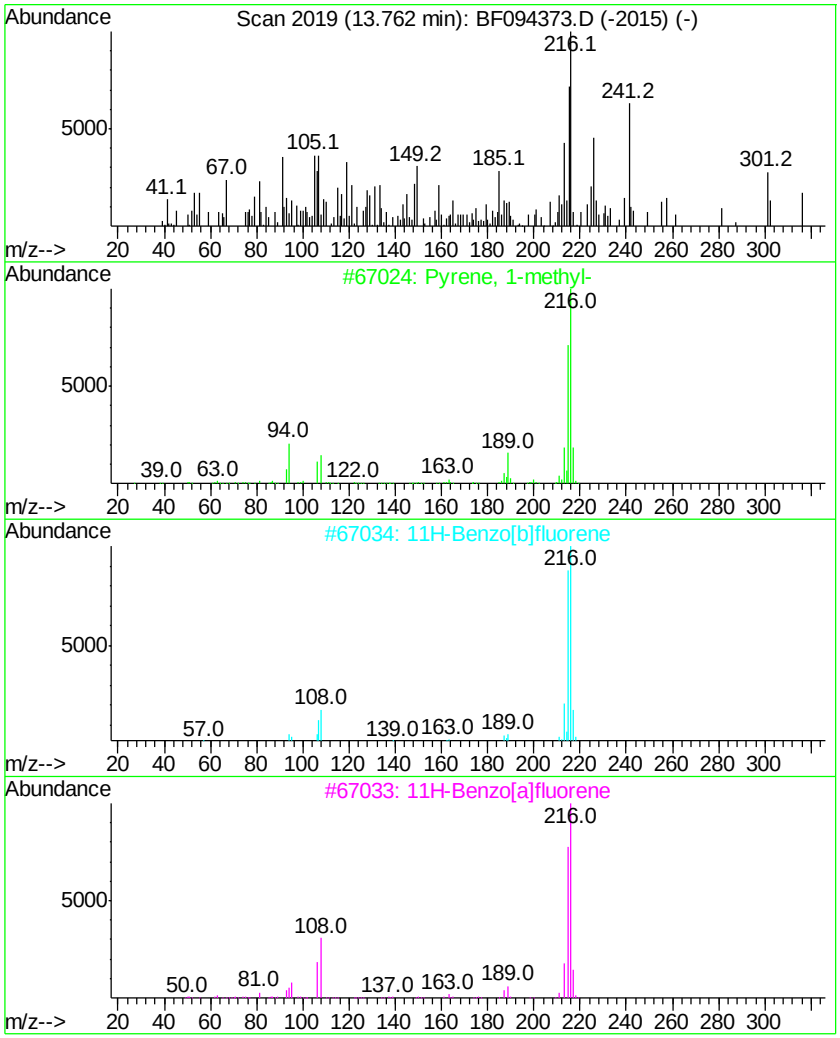
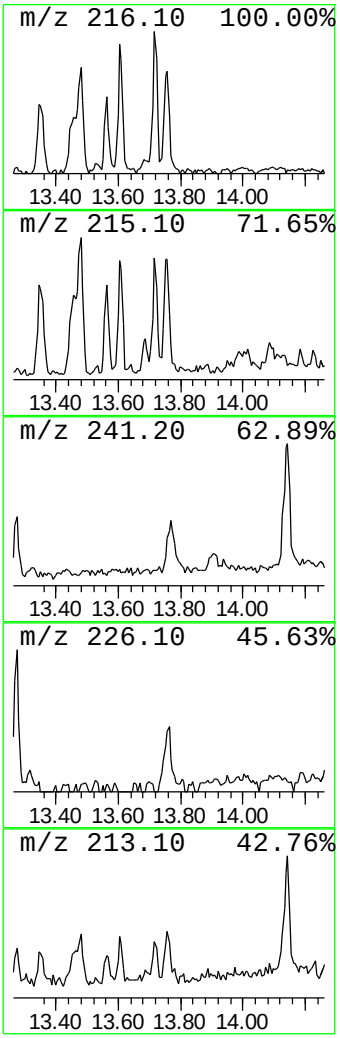
Peak Number 10 Pyrene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	2.80 ng	144799	Chrysene-d12	14.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	83
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	58
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	58
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	52
5	1,3,5-Triazine-2,4,6-triamine, N...	216	C10H12N6	046731-79-7	50





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Acq On : 12 Apr 2017 1:50
Operator : SJ/MA
Sample : I2610-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SOURCE3-SE

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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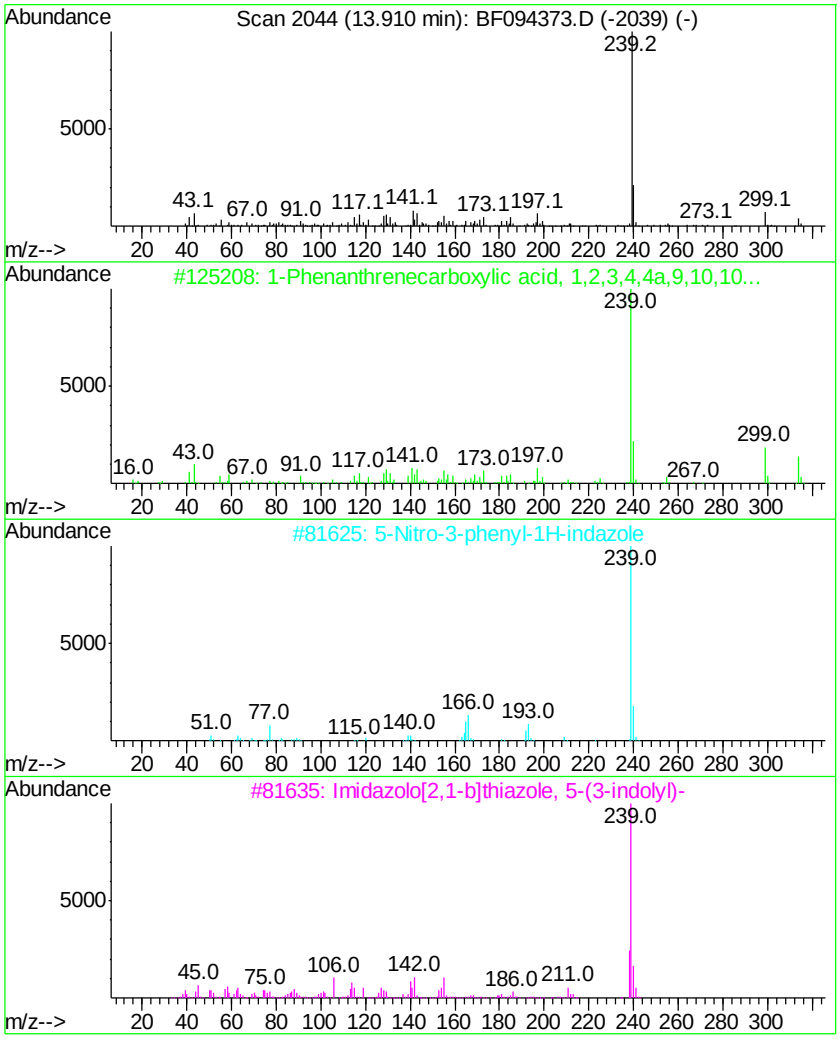
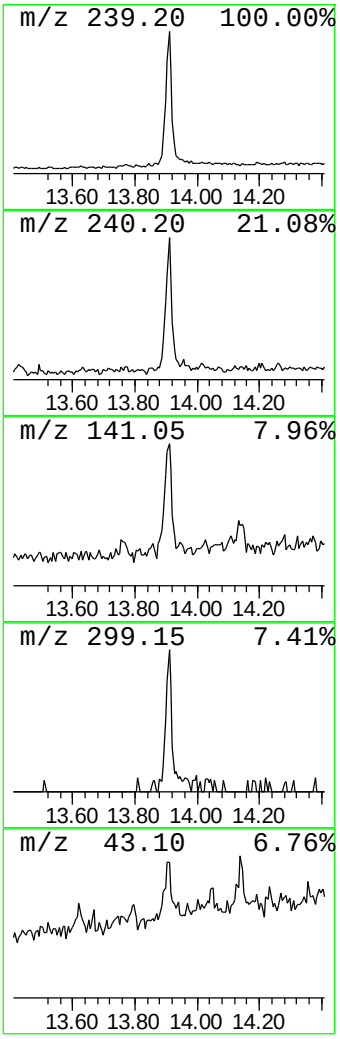
Peak Number 11 1-Phenanthrenecarboxylic ac... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.91	5.21 ng	269629	Chrysene-d12	14.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenanthrenecarboxylic acid, 1...	314	C21H30O2	001235-74-1	95
2	5-Nitro-3-phenyl-1H-indazole	239	C13H9N3O2	293758-67-5	64
3	Imidazolo[2,1-b]thiazole, 5-(3-i...	239	C13H9N3S	327097-37-0	59
4	Benzaldehyde, 3-[4-(1,1-dimethyl...	254	C17H18O2	069770-23-6	50
5	4-(4-Oxo-1,2,3,4,6,7,12,12b-octa...	326	C19H22N2O3	1000137-91-1	45





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Sample : I2610-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SOURCE3-SE

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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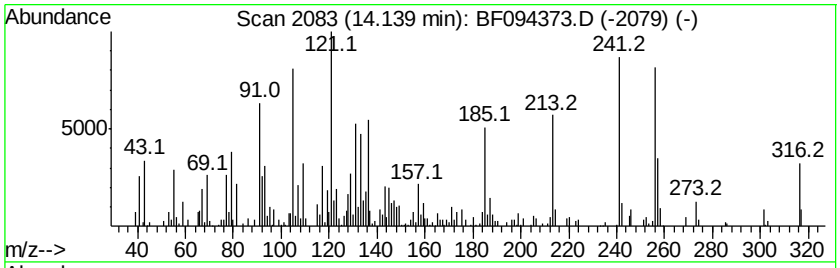
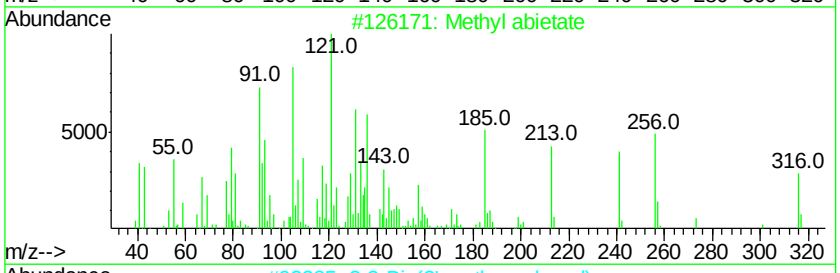
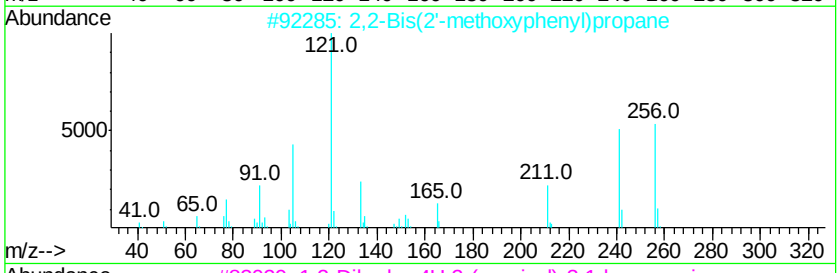
Peak Number 12 Methyl abietate Concentration Rank 5

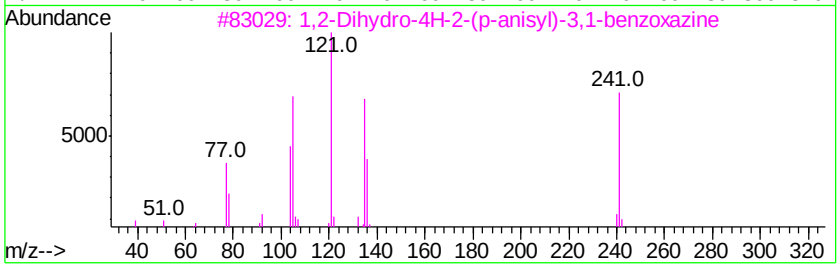
R.T.	EstConc	Area	Relative to ISTD	R.T.
14.14	4.72 ng	244491	Chrysene-d12	14.55

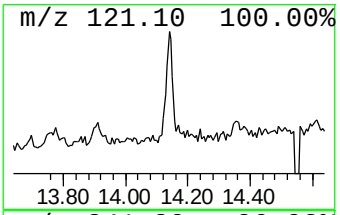
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl abietate	316	C21H32O2	000127-25-3	90
2	2,2-Bis(2'-methoxyphenyl)propane	256	C17H20O2	056751-16-7	43
3	1,2-Dihydro-4H-2-(p-anisyl)-3,1-...	241	C15H15NO2	082085-87-8	30
4	Diazene, (4-methoxy-2-methylphen...	256	C15H16N2O2	029418-53-9	30
5	2,4,6-Octatriene, 2,6-dimethyl-,...	136	C10H16	007216-56-0	25

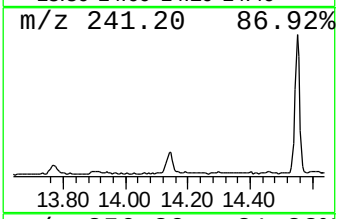


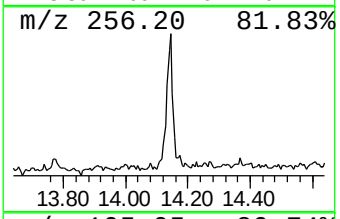


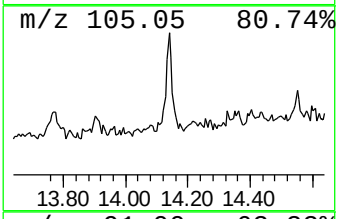


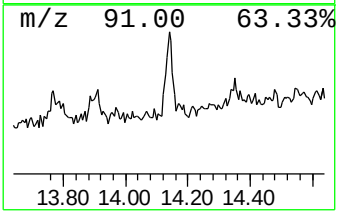












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Sample : I2610-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SOURCE3-SE

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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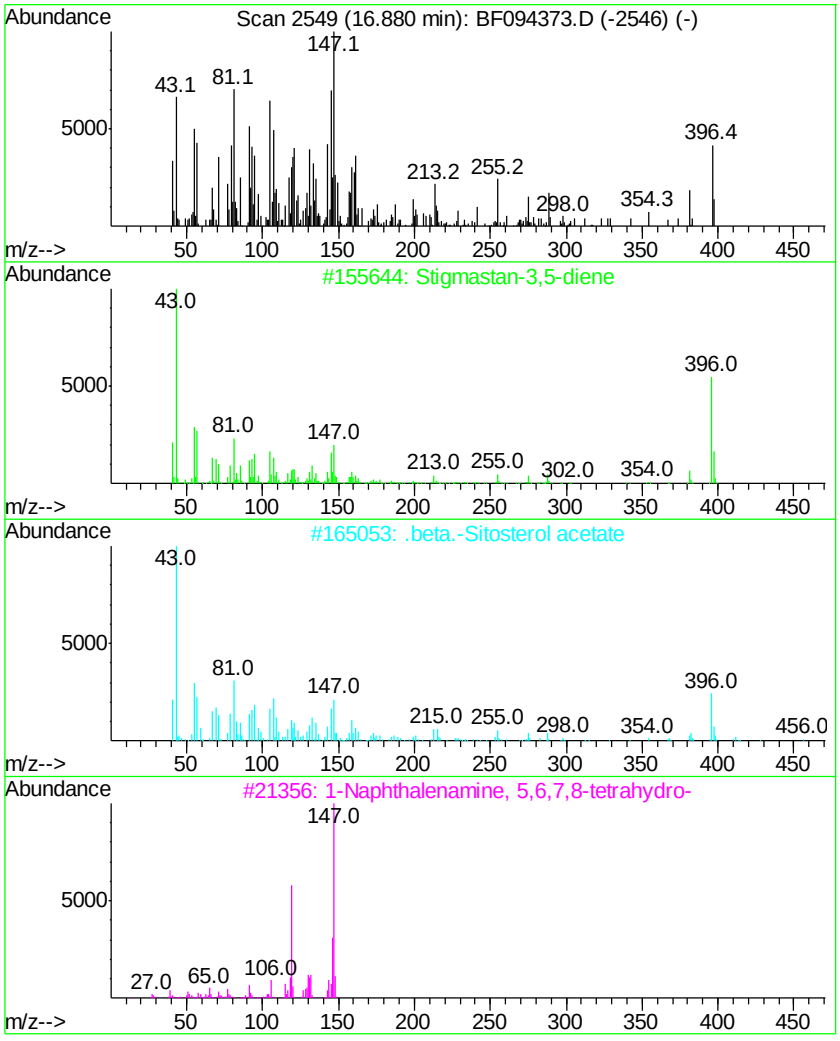
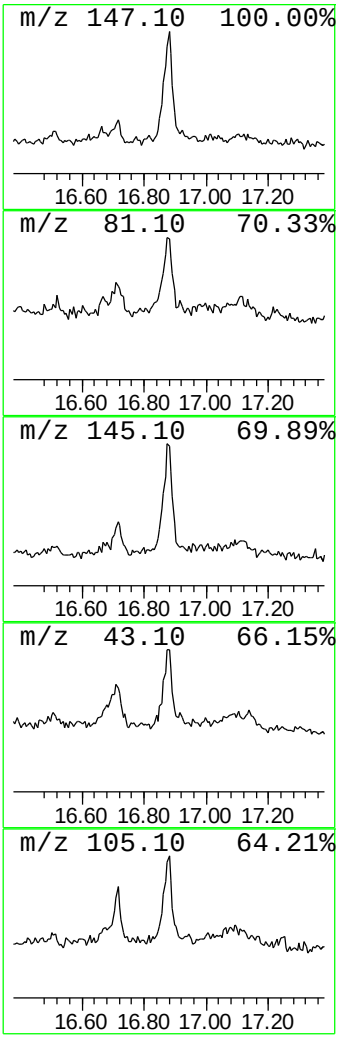
Peak Number 13 Stigmastan-3,5-diene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.88	15.93 ng	527310	Perylene-d12	16.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Stigmastan-3,5-diene	396	C29H48	1000214-16-4	89
2	.beta.-Sitosterol acetate	456	C31H52O2	000915-05-9	25
3	1-Naphthalenamine, 5,6,7,8-tetra...	147	C10H13N	002217-41-6	18
4	Acetic acid 2-oxo-2,3-dihydro-1H...	218	C11H10N2O3	016780-58-8	12
5	Benzenesulfonic acid, 4-nitro-, ...	429	C17H20BrNO5S	274255-39-9	10





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
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Benzenesulfonamid...	10.60	2.6	ng	126503	4	11.30	958713	20.0
4H-Cyclopenta[def...	12.05	2.8	ng	132548	4	11.30	958713	20.0
Phosphonic acid, ...	12.49	2.1	ng	98965	4	11.30	958713	20.0
Bis(acetylacetona...	12.64	4.4	ng	210638	4	11.30	958713	20.0
Borane, 2,3-dimet...	12.72	3.0	ng	146011	4	11.30	958713	20.0
Cyclohexadecane	12.93	2.6	ng	136954	5	14.55	1035990	20.0
Pyrene, 1-methyl-	13.76	2.8	ng	144799	5	14.55	1035990	20.0
1-Phenanthrenecar...	13.91	5.2	ng	269629	5	14.55	1035990	20.0
Methyl abietate	14.14	4.7	ng	244491	5	14.55	1035990	20.0
Stigmastan-3,5-diene	16.88	15.9	ng	527310	6	16.18	661981	20.0