LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052319\
Data File : BF114554.D
Acq On : 24 May 2019 5:04
Operator : JU/SJ
Sample : K3008-01 2X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : COMP-19

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-BF051019.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	5.439	567	570	589	rBV	692622	921122	28.31%	2.403%
2	6.457	740	743	762	rBV	1012105	1134906	34.88%	2.961%
3	6.586	762	765	774	rVV	1168377	1067325	32.80%	2.785%
4	6.810	800	803	811	rBV	1090472	1024353	31.48%	2.673%
5	6.969	826	830	839	rBV	967480	843409	25.92%	2.200%
6	7.375	896	899	911	rBV	664274	655816	20.16%	1.711%
7	8.092	1018	1021	1031	rBV	1565185	1370291	42.12%	3.575%
8	9.169	1200	1204	1210	rBV	1493595	1229707	37.79%	3.208%
9	9.563	1268	1271	1280	rVB	151972	141810	4.36%	0.370%
10	9.851	1316	1320	1323	rBV	1829237	1503134	46.20%	3.922%
11	9.880	1323	1325	1334	rVB	131541	120089	3.69%	0.313%
12	10.057	1352	1355	1359	rVB	40681	34645	1.06%	0.090%
13	10.398	1410	1413	1418	rVB	97153	94642	2.91%	0.247%
14	10.639	1451	1454	1465	rVB	549851	561245	17.25%	1.464%
15	10.769	1473	1476	1481	rVB3	78334	88217	2.71%	0.230%
16	10.821	1481	1485	1487	rBV	123609	117970	3.63%	0.308%
17	10.851	1487	1490	1494	rVV2	127331	173407	5.33%	0.452%
18	10.886	1494	1496	1498	rVV	151287	129116	3.97%	0.337%
19	10.910	1498	1500	1503	rVV	89502	86739	2.67%	0.226%
20	10.951	1503	1507	1509	rVV3	73942	114317	3.51%	0.298%
21	10.998	1512	1515	1519	rVV3	126009	161111	4.95%	0.420%
22	11.033	1519	1521	1524	rVV	152628	147247	4.53%	0.384%
23	11.074	1526	1528	1537	rVV2	106665	134518	4.13%	0.351%
24	11.151	1538	1541	1545	rVV	52287	48618	1.49%	0.127%
25	11.192	1545	1548	1551	rVV4	53460	69726	2.14%	0.182%
26	11.233	1552	1555	1559	rVB	75999	82776	2.54%	0.216%
27	11.339	1569	1573	1575	rBV	1758070	1551771	47.69%	4.049%
28	11.363	1575	1577	1581	rVV	1466854	1145006	35.19%	2.987%
29	11.416	1583	1586	1589	rVV	327627	288433	8.86%	0.753%
30	11.451	1589	1592	1598	rVB2	22881	41516	1.28%	0.108%
31	11.574	1608	1613	1619	rBV2	151080	190881	5.87%	0.498%
32	11.627	1619	1622	1624	rVV2	45844	49761	1.53%	0.130%
33	11.680	1627	1631	1635	rVB3	31035	52440	1.61%	0.137%
34	11.839	1655	1658	1660	rBV	182480	147788	4.54%	0.386%

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Misc :
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Instrument : BNA_F
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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 >
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Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.868	1660	1663	1668	rVB	271428	242703	7.46%	0.633%
36	11.915	1668	1671	1674	rBV	91281	77568	2.38%	0.202%
37	11.951	1674	1677	1680	rBV	335824	354788	10.90%	0.926%
38	12.027	1686	1690	1692	rBV3	38201	48611	1.49%	0.127%
39	12.133	1705	1708	1711	rBV	140055	114863	3.53%	0.300%
40	12.174	1712	1715	1718	rVB	90679	93929	2.89%	0.245%
41	12.210	1718	1721	1724	rBV2	29833	33389	1.03%	0.087%
42	12.280	1731	1733	1737	rVB	47286	42200	1.30%	0.110%
43	12.315	1737	1739	1741	rBV	57236	45789	1.41%	0.119%
44	12.339	1741	1743	1748	rBV2	37610	36386	1.12%	0.095%
45	12.392	1749	1752	1754	rBV	89881	108188	3.33%	0.282%
46	12.427	1755	1758	1760	rVV3	107004	120703	3.71%	0.315%
47	12.486	1764	1768	1771	rBV2	123769	148401	4.56%	0.387%
48	12.557	1776	1780	1783	rBV	3142357	2794786	85.90%	7.292%
49	12.621	1789	1791	1793	rVB	55650	41341	1.27%	0.108%
50	12.704	1803	1805	1808	rVB	85713	68820	2.12%	0.180%
51	12.786	1813	1819	1822	rBV	2637023	3066585	94.25%	8.001%
52	12.833	1825	1827	1832	rVB2	130618	124703	3.83%	0.325%
53	12.915	1836	1841	1844	rBV2	1175546	1136070	34.92%	2.964%
54	12.945	1844	1846	1850	rVB	147448	150652	4.63%	0.393%
55	12.998	1851	1855	1859	rBV2	171219	289376	8.89%	0.755%
56	13.110	1867	1874	1877	rBV2	421798	663690	20.40%	1.732%
57	13.174	1883	1885	1888	rBV	250947	215133	6.61%	0.561%
58	13.215	1888	1892	1899	rBV4	308856	576136	17.71%	1.503%
59	13.310	1905	1908	1911	rBV	143004	170558	5.24%	0.445%
60	13.339	1911	1913	1916	rBV2	140245	140124	4.31%	0.366%
61	13.533	1944	1946	1949	rVB	131888	107259	3.30%	0.280%
62	13.627	1959	1962	1966	rBV	286790	254736	7.83%	0.665%
63	13.733	1977	1980	1982	rBV	370106	336893	10.35%	0.879%
64	13.757	1982	1984	1986	rVV	267701	219555	6.75%	0.573%
65	13.786	1986	1989	1995	rVV3	240726	460802	14.16%	1.202%
66	13.839	1995	1998	2004	rVB	358571	360209	11.07%	0.940%
67	13.980	2015	2022	2025	rBV2	2024969	3253662	100.00%	8.489%
68	14.009	2025	2027	2031	rVB	1393016	1299225	39.93%	3.390%
69	14.092	2038	2041	2044	rVB	349900	272530	8.38%	0.711%
70	14.209	2058	2061	2065	rBV2	158956	226534	6.96%	0.591%
71	14.374	2087	2089	2091	rVB	256683	190760	5.86%	0.498%

LSC Area Percent Report

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Misc :
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Instrument : BNA_F
ClientSampleId : COMP-19

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-BF051019.M
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72	15.033	2197	2201	2204	rBV	1059971	1586189	48.75%	4.138%
73	15.186	2224	2227	2230	rVB	395265	387230	11.90%	1.010%
74	15.333	2248	2252	2258	rBV	571433	679929	20.90%	1.774%
75	15.398	2259	2263	2267	rVB	737176	933998	28.71%	2.437%
76	15.456	2269	2273	2276	rVV	573456	711581	21.87%	1.857%
77	16.939	2520	2525	2533	rVB2	324238	618450	19.01%	1.614%

Sum of corrected areas: 38328886

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052319\
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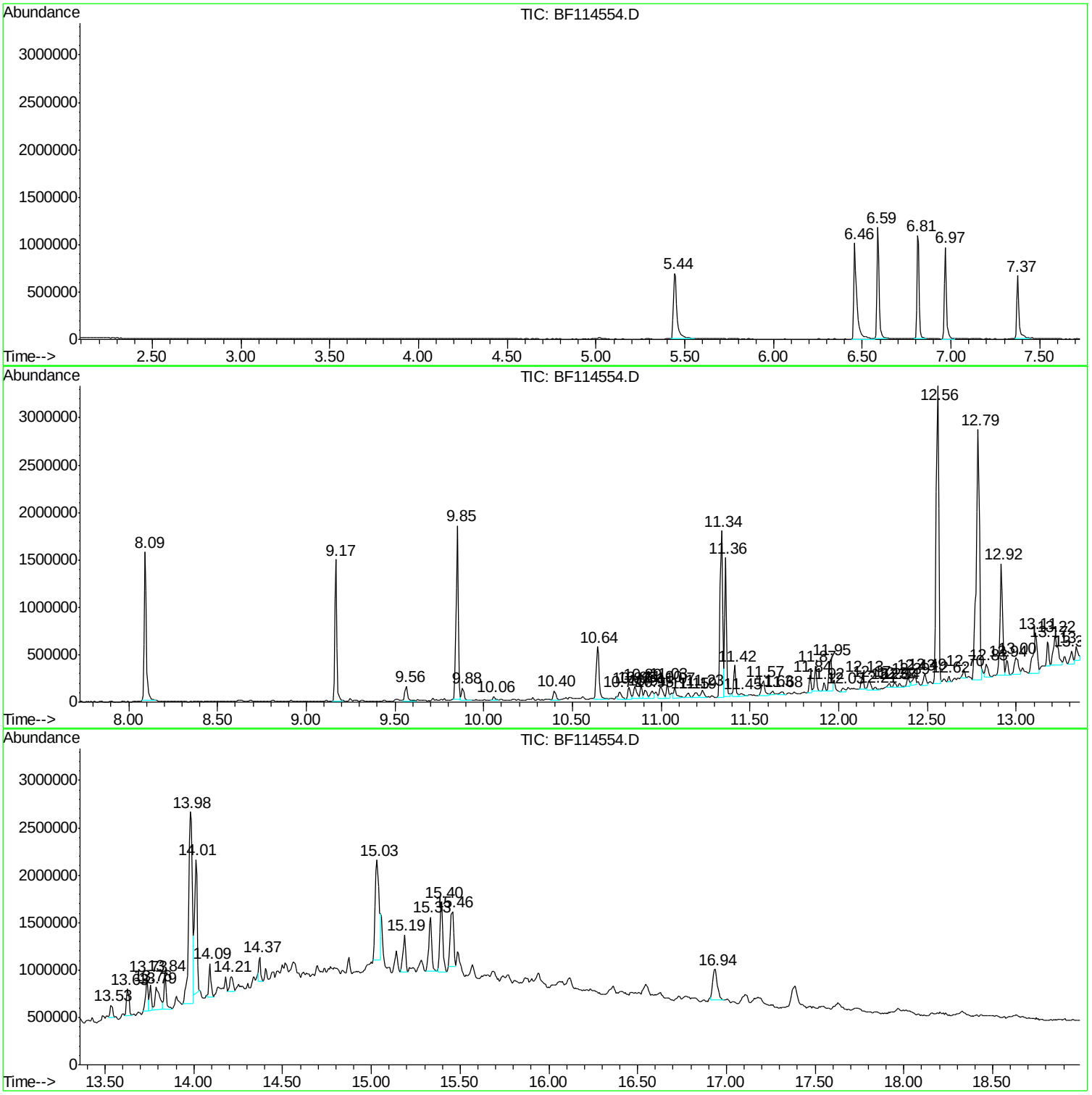
Sample : K3008-01 2X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMP-19

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.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\BF052319\
Data File : BF114554.D
Acq On : 24 May 2019 5:04
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Misc :
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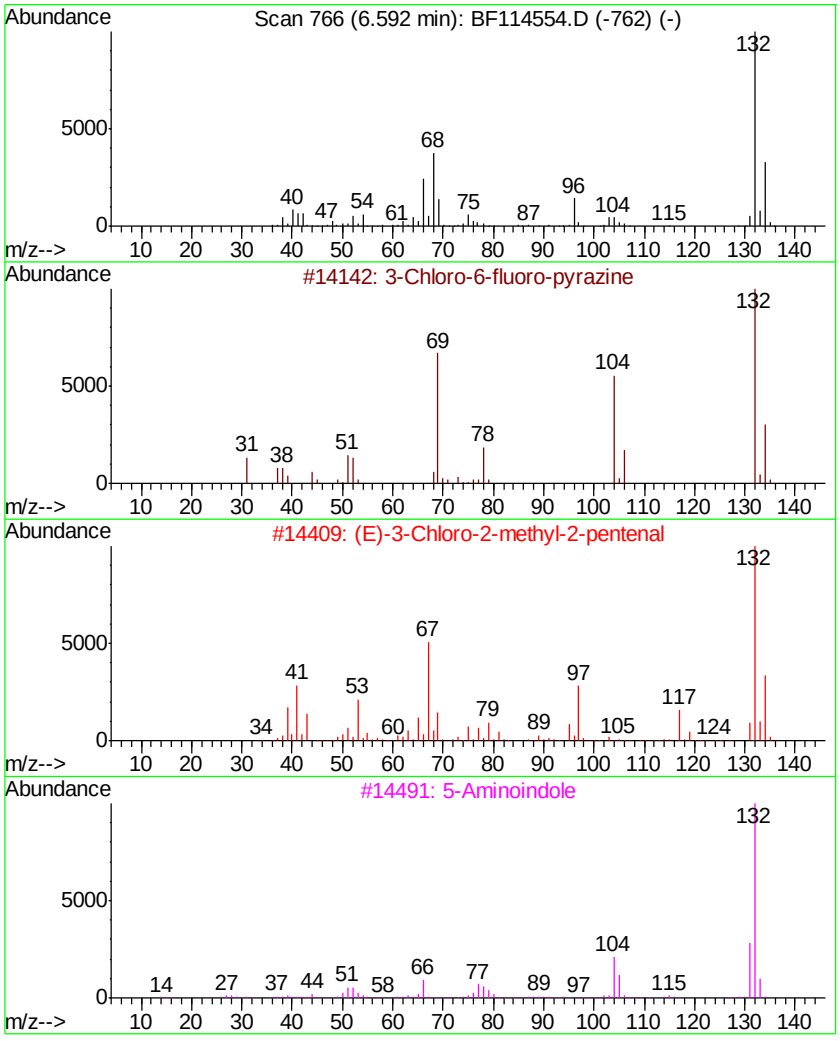
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

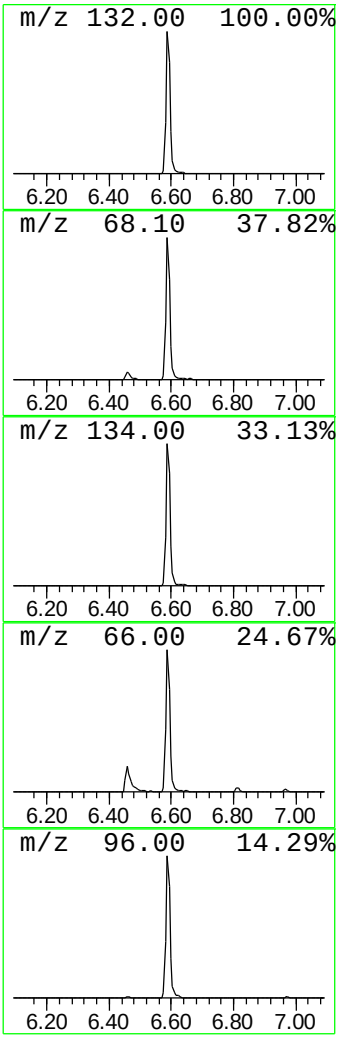
****

Peak Number 1 unknown6.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	20.84 ng	1067330	1,4-Dichlorobenzene-d4	6.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3	5-Aminoindole	132	C8H8N2	005192-03-0	12
4	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5	1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052319\
Data File : BF114554.D
Acq On : 24 May 2019 5:04
Operator : JU/SJ
Sample : K3008-01 2X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : COMP-19

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

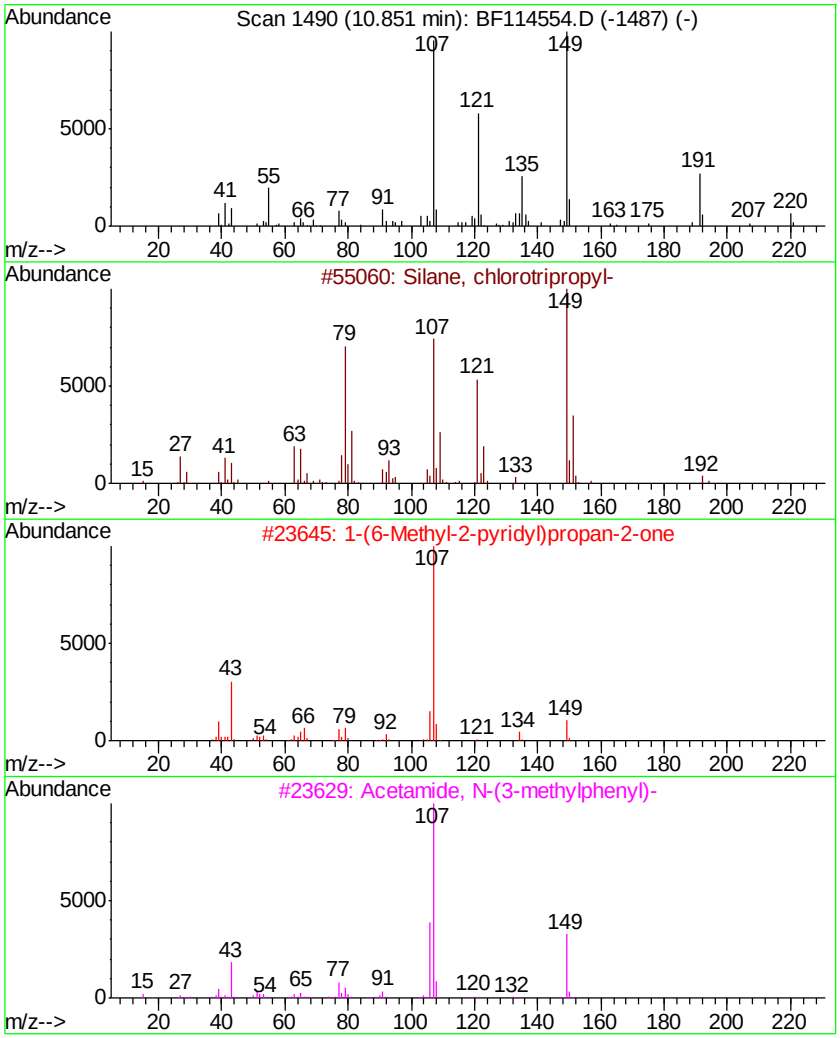
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

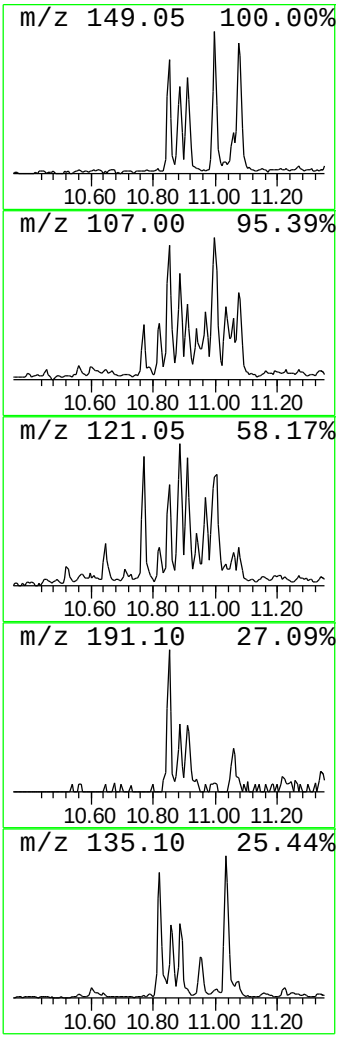
****

Peak Number 3 Silane, chlorotripropyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.85	2.23 ng	173407	Phenanthrene-d10	11.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Silane, chlorotripropyl-	192	C9H21ClSi	000995-25-5	64
2	1-(6-Methyl-2-pyridyl)propan-2-one	149	C9H11NO	065702-08-1	58
3	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
4	5H-Pyrrolo(3,2-d)pyrimidine-2,4-...	149	C6H7N5	1000244-21-4	53
5	1,3-Cyclopentadiene, 5,5-dimethy...	178	C13H22	1000163-88-0	38





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052319\
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ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_F
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
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TIC Library : C:\DATABASE\NIST11.L
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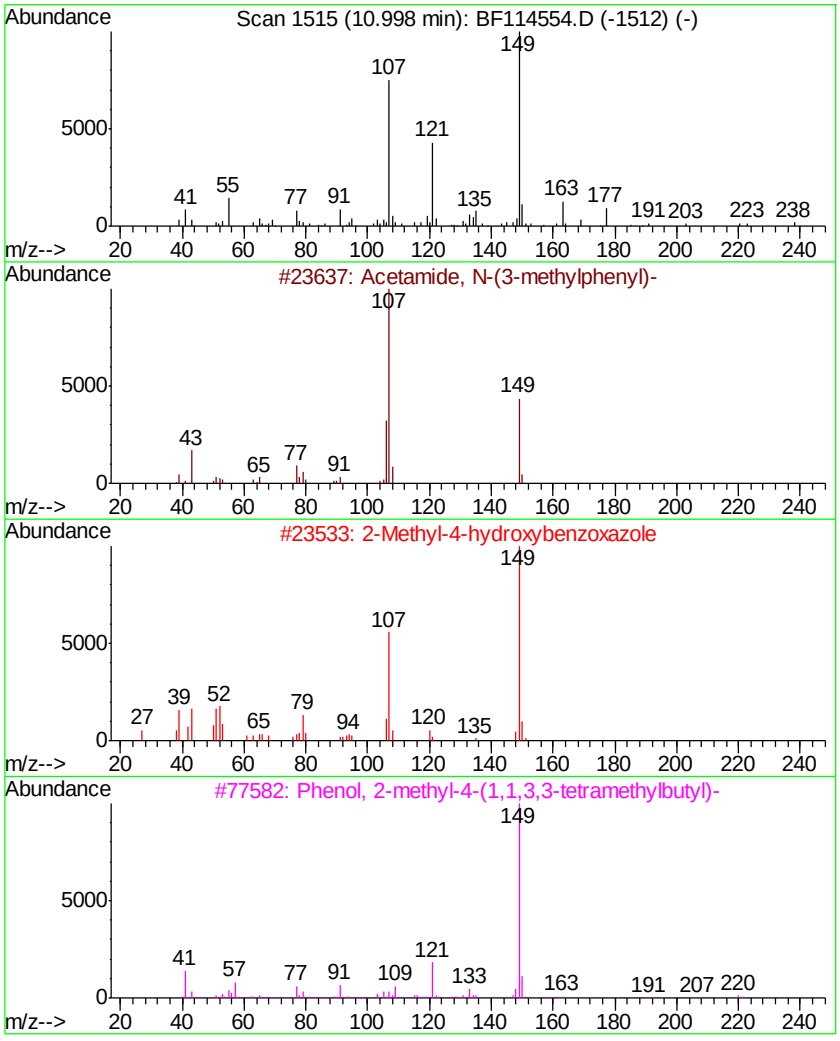
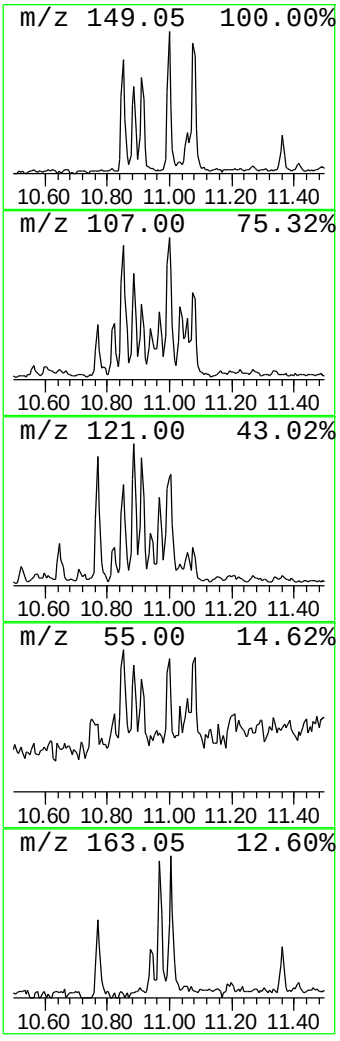
****

Peak Number 4 Acetamide, N-(3-methylphenyl)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.00	2.08 ng	161111	Phenanthrene-d10	11.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	64
2	2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	47
3	Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	43
4	2H-Indol-2-one, 1,3-dihydro-5-hy...	149	C8H7NO2	003416-18-0	38
5	1,3-Cyclopentadiene, 5,5-dimethy...	178	C13H22	1000163-88-0	38





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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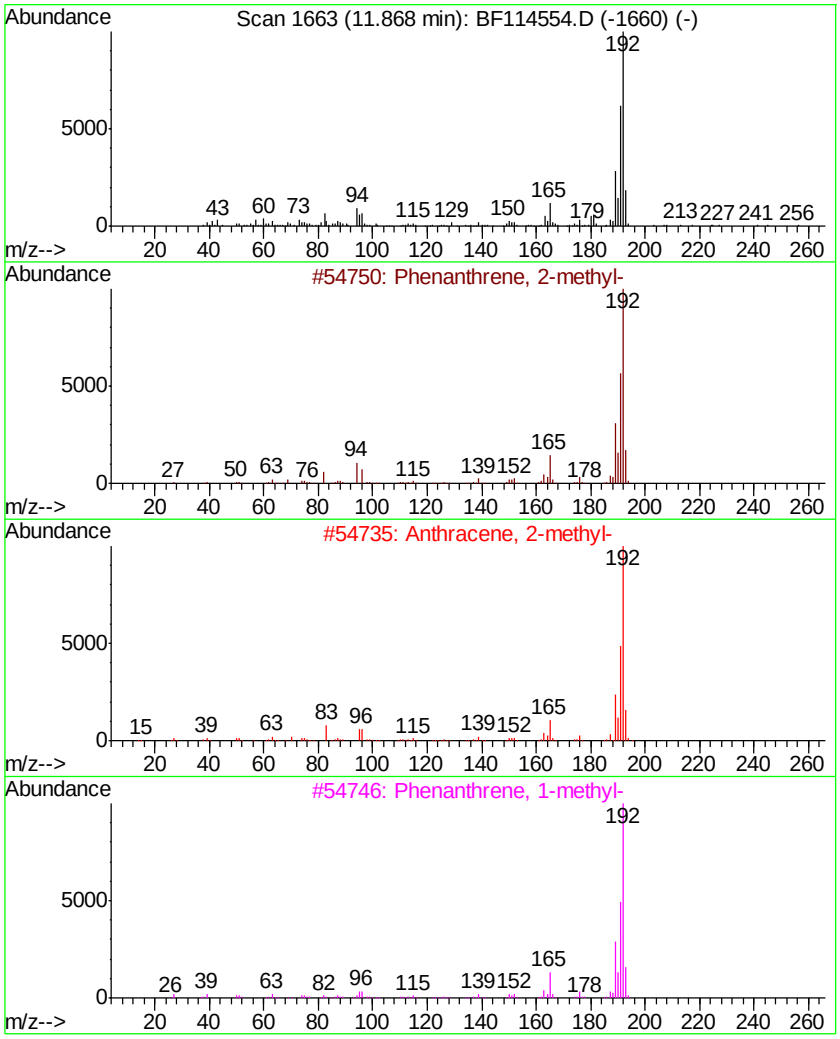
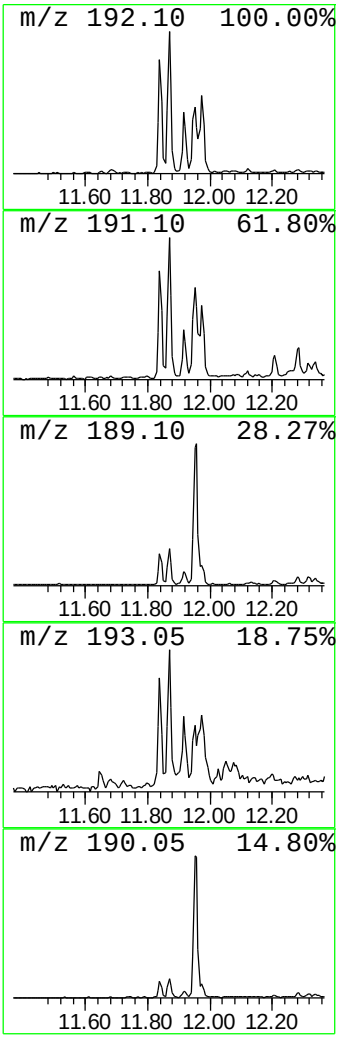
Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	3.13 ng	242703	Phenanthrene-d10	11.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	97
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	96
5	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	95





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Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : COMP-19

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

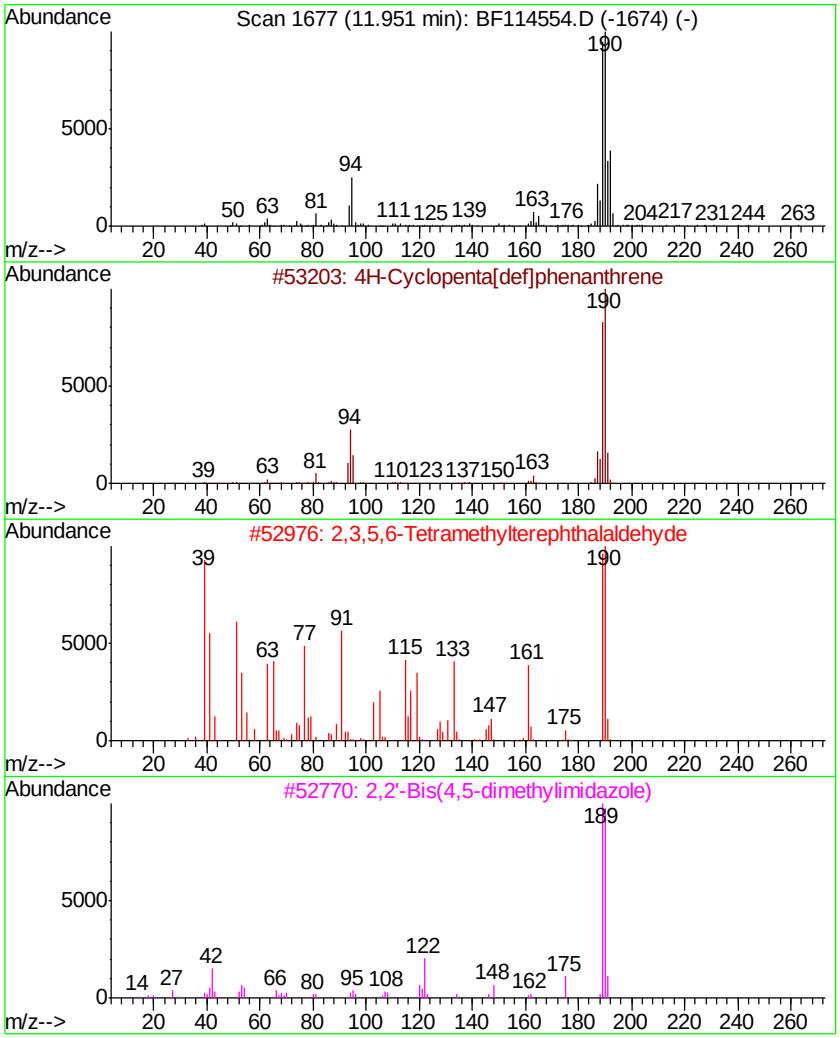
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

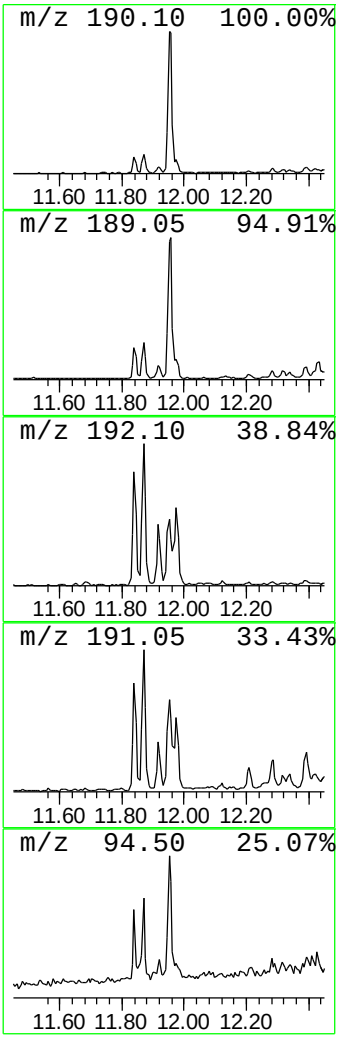
****

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.95	4.57 ng	354788	Phenanthrene-d10	11.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2	2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	47
3	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	38
4	1-(Phenylethynyl)-1H-1,2,3-benzo...	219	C14H9N3	209912-23-2	22
5	Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	17





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Data File : BF114554.D
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Sample : K3008-01 2X
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Instrument : BNA_F
ClientSampleId : COMP-19

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

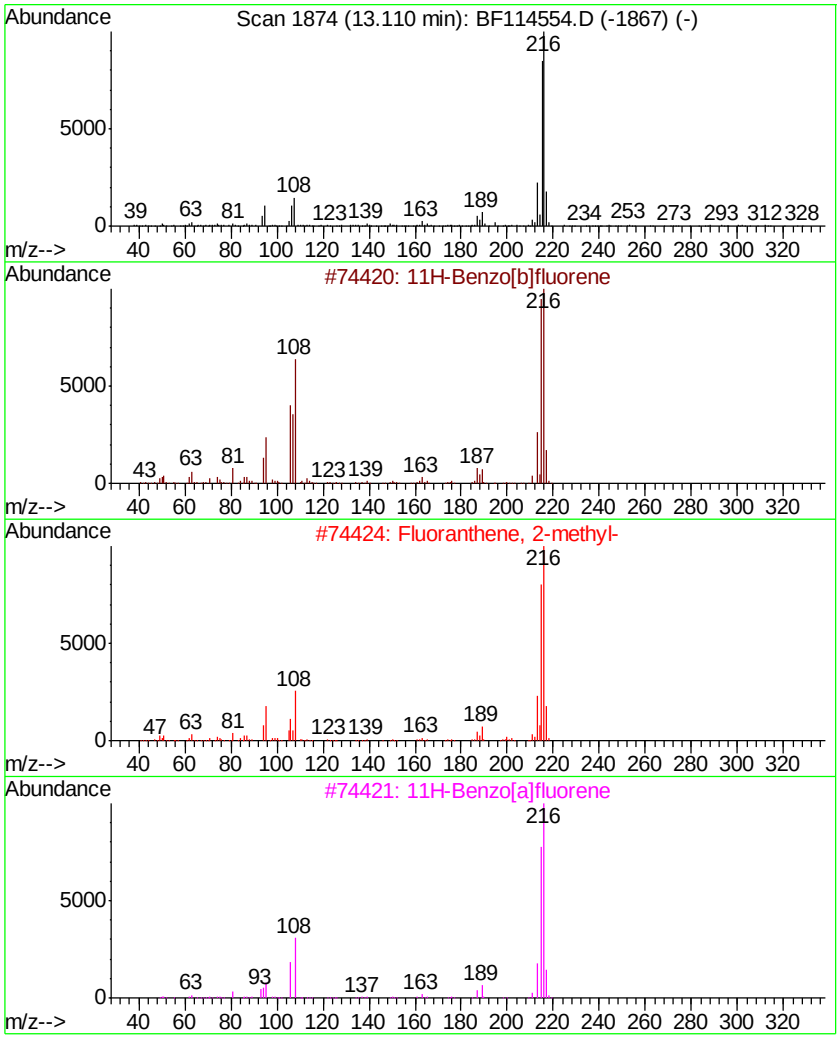
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

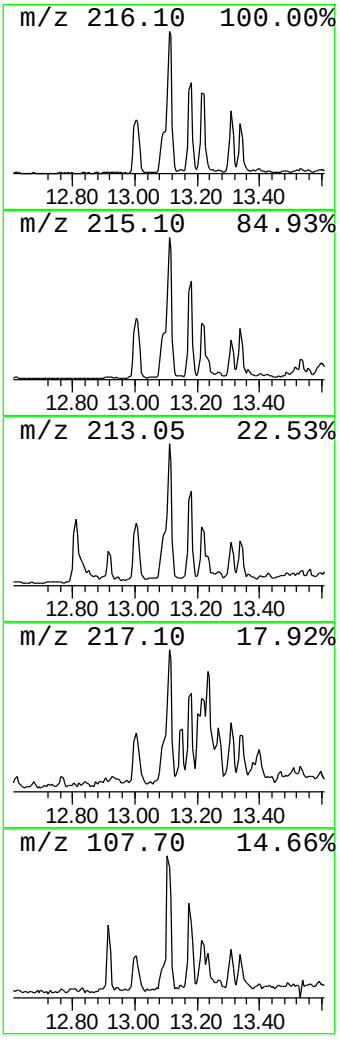
****

Peak Number 7 11H-Benzo[b]fluorene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.11	4.08 ng	663690	Chrysene-d12	13.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
4	Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
5	Pyrene, 4-methyl-	216	C17H12	003353-12-6	60





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052319\
Data File : BF114554.D
Acq On : 24 May 2019 5:04
Operator : JU/SJ
Sample : K3008-01 2X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : COMP-19

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

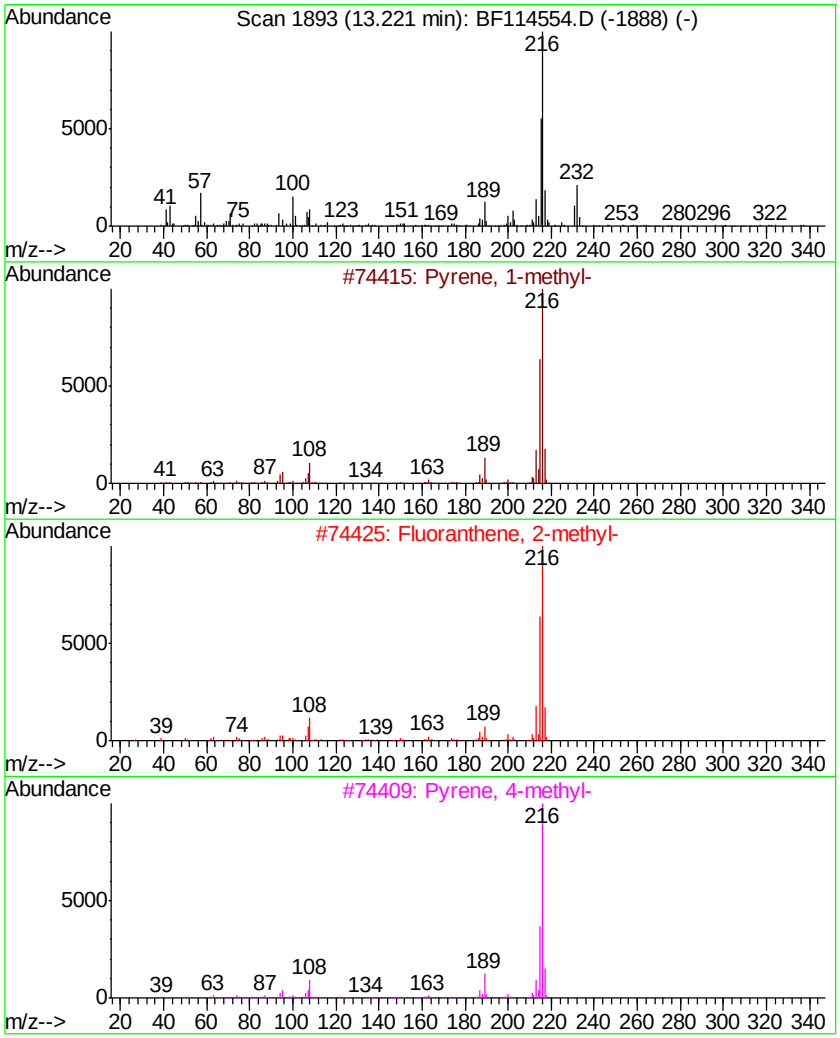
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

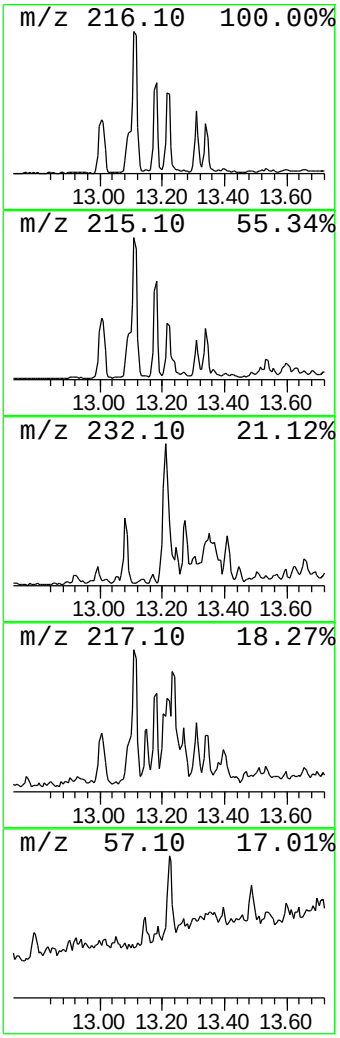
****

Peak Number 8 Pyrene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	3.54 ng	576136	Chrysene-d12	13.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	95
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
3	Pyrene, 4-methyl-	216	C17H12	003353-12-6	86
4	Pyrene, 2-methyl-	216	C17H12	003442-78-2	86
5	: 4-((2-Methylphenyl)diazenyl)-1...	216	C10H12N6	1000332-58-9	74





Library Search Compound Report

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Data File : BF114554.D
Acq On : 24 May 2019 5:04
Operator : JU/SJ
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Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : COMP-19

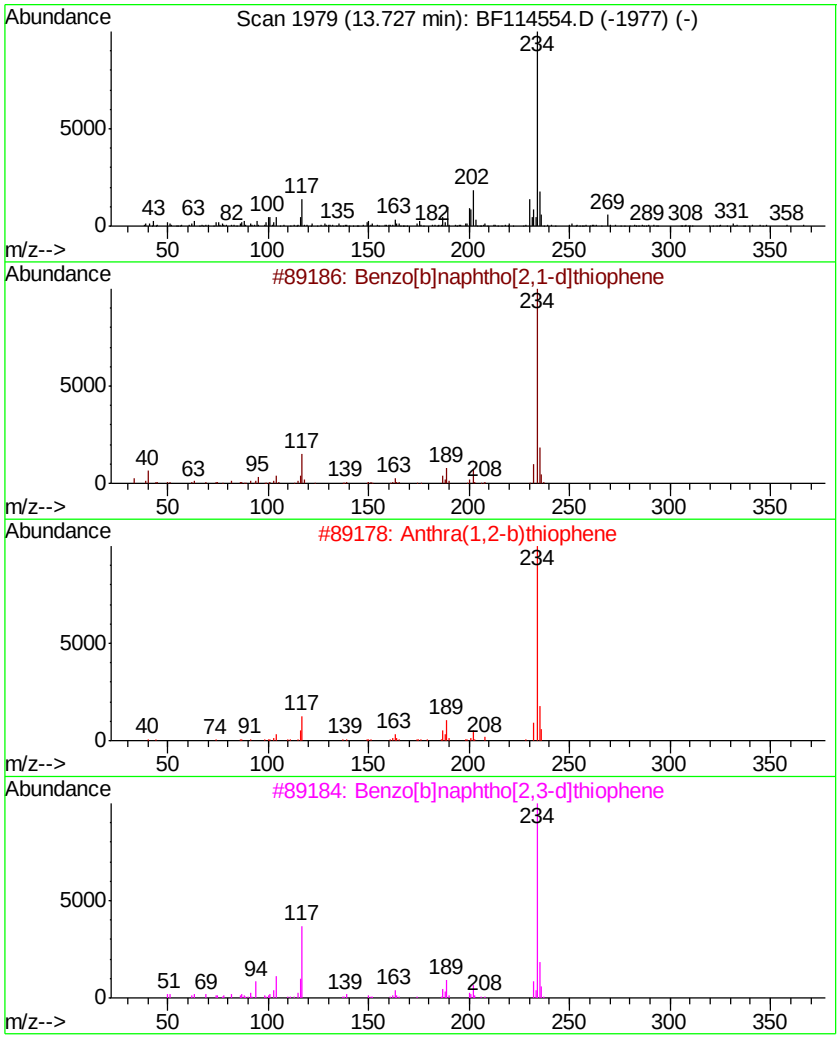
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

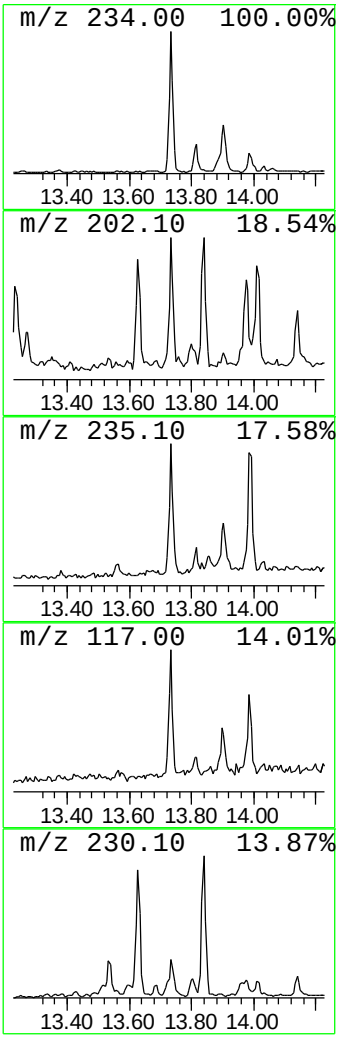
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	2.07 ng	336893	Chrysene-d12	13.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	95
2	Anthra(1,2-b)thiophene	234	C16H10S	000227-86-1	58
3	Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	58
4	Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	58
5	Quinoxalino[2,3-b]quinoxaline, 5...	234	C14H10N4	000531-46-4	47





Library Search Compound Report

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Data File : BF114554.D
Acq On : 24 May 2019 5:04
Operator : JU/SJ
Sample : K3008-01 2X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : COMP-19

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
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TIC Library : C:\DATABASE\NIST11.L
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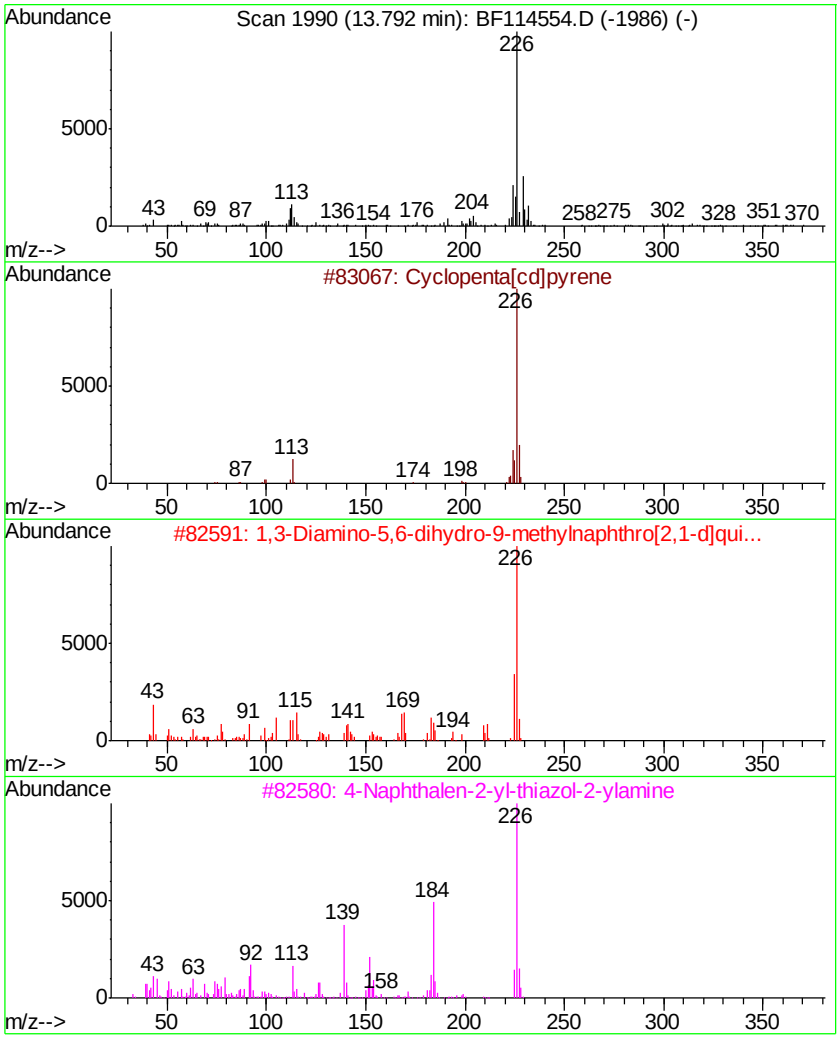
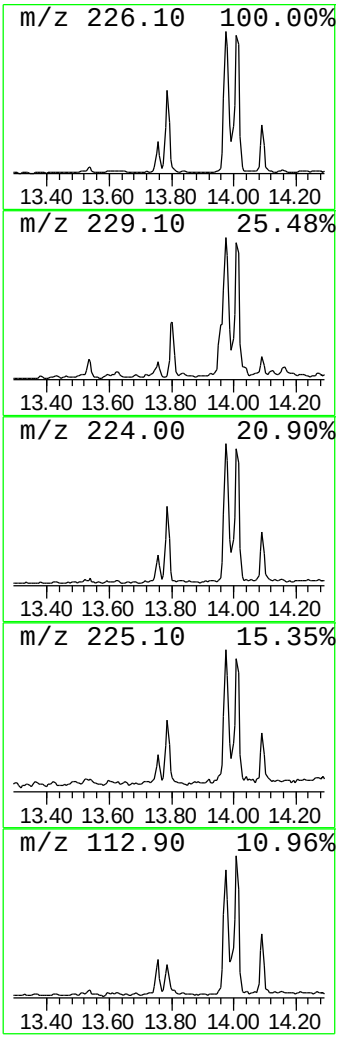
Peak Number 10 Cyclopenta[cd]pyrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	2.83 ng	460802	Chrysene-d12	13.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	70
2	1,3-Diamino-5,6-dihydro-9-methyl...	226	C13H14N4	037436-39-8	42
3	4-Naphthalen-2-yl-thiazol-2-ylamine	226	C13H10N2S	1000300-53-4	40
4	1,8-Diazacyclotetradecane-2,7-dione	226	C12H22N2O2	004266-66-4	37
5	.beta.-Carboline, 7-methoxy-1,2-...	226	C14H14N2O	006519-18-2	33





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Acq On : 24 May 2019 5:04
Operator : JU/SJ
Sample : K3008-01 2X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : COMP-19

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
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TIC Library : C:\DATABASE\NIST11.L
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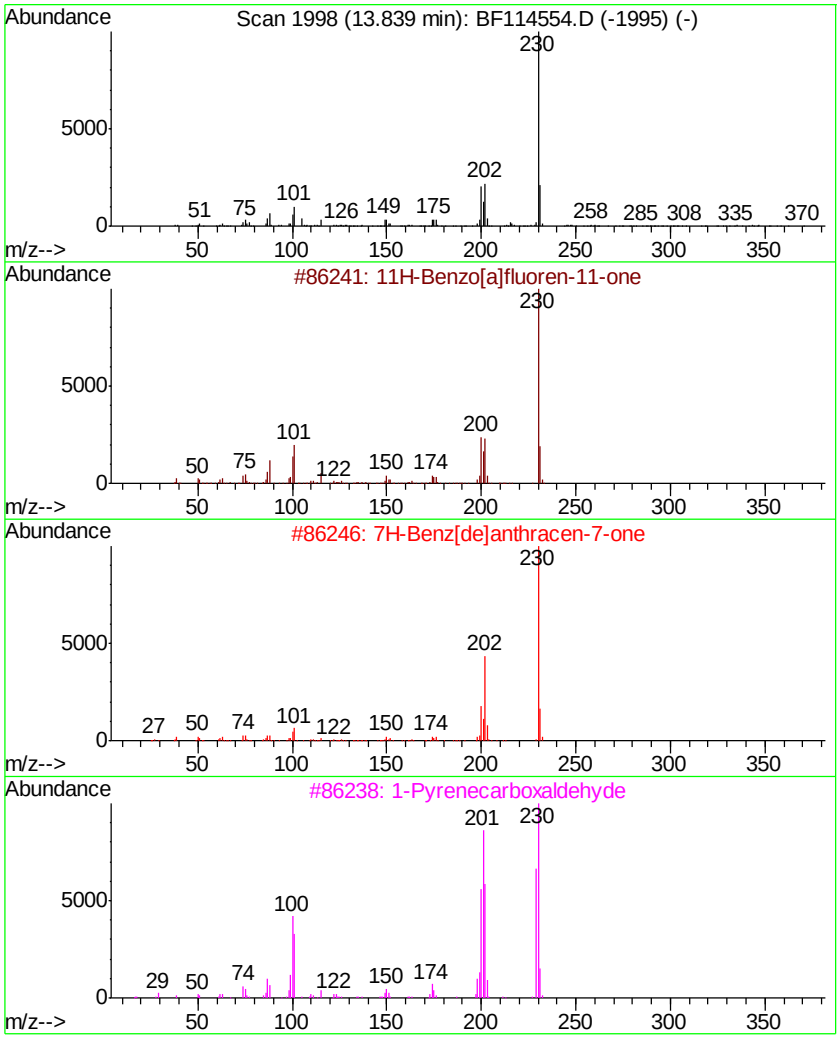
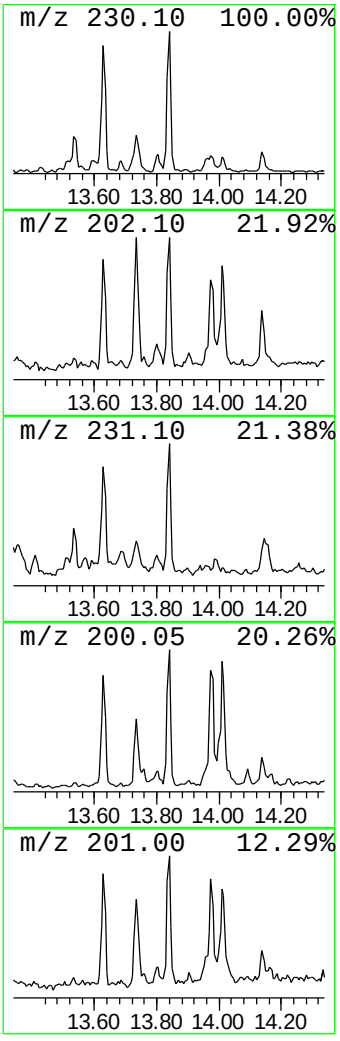
Peak Number 11 11H-Benzo[a]fluoren-11-one Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.84	2.21 ng	360209	Chrysene-d12	13.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	93
2	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
3	1-Pyrenecarboxaldehyde	230	C17H10O	003029-19-4	78
4	p-Terphenyl	230	C18H14	000092-94-4	53
5	8-Chlorobenzo[h][1,6]-naphthyrid...	230	C12H7ClN2O	025100-26-9	53





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052319\
Data File : BF114554.D
Acq On : 24 May 2019 5:04
Operator : JU/SJ
Sample : K3008-01 2X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : COMP-19

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.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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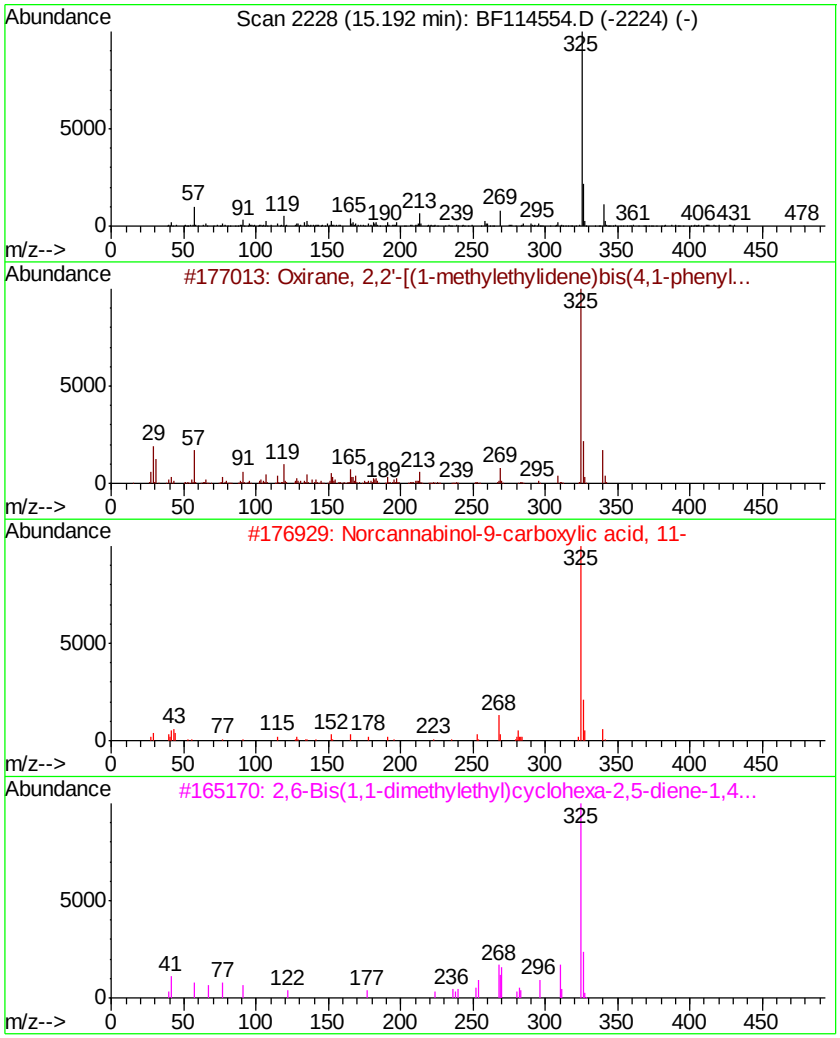
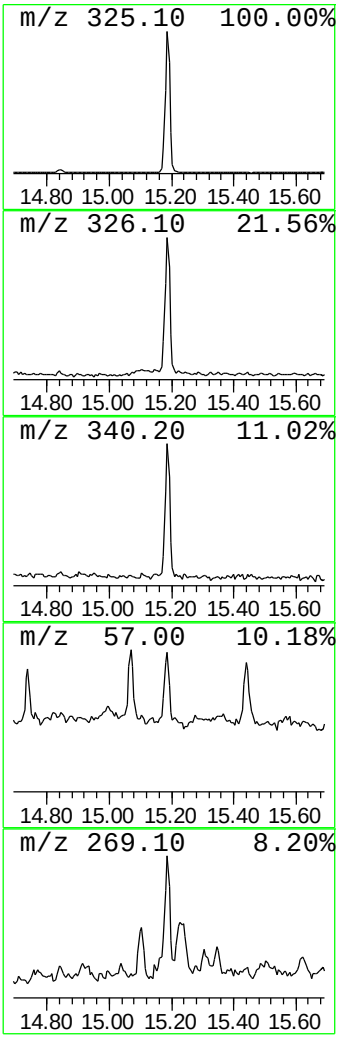
Peak Number 12 Oxirane, 2,2'-[(1-methyleth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	10.88 ng	387230	Perylene-d12	15.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, 2,2'-[(1-methylethylide...	340	C21H24O4	001675-54-3	96
2	Norcannabinol-9-carboxylic acid,...	340	C21H24O4	053989-32-5	64
3	2,6-Bis(1,1-dimethylethyl)cycloh...	325	C21H27NO2	017119-01-6	64
4	Butane, 1,4-bis(dicyclopentylpho...	394	C24H44P2	163106-88-5	59
5	Bis(pentamethylcyclopentadienyl)...	325	C20H30Mn	067506-86-9	59





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Operator : JU/SJ
Sample : K3008-01 2X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : COMP-19

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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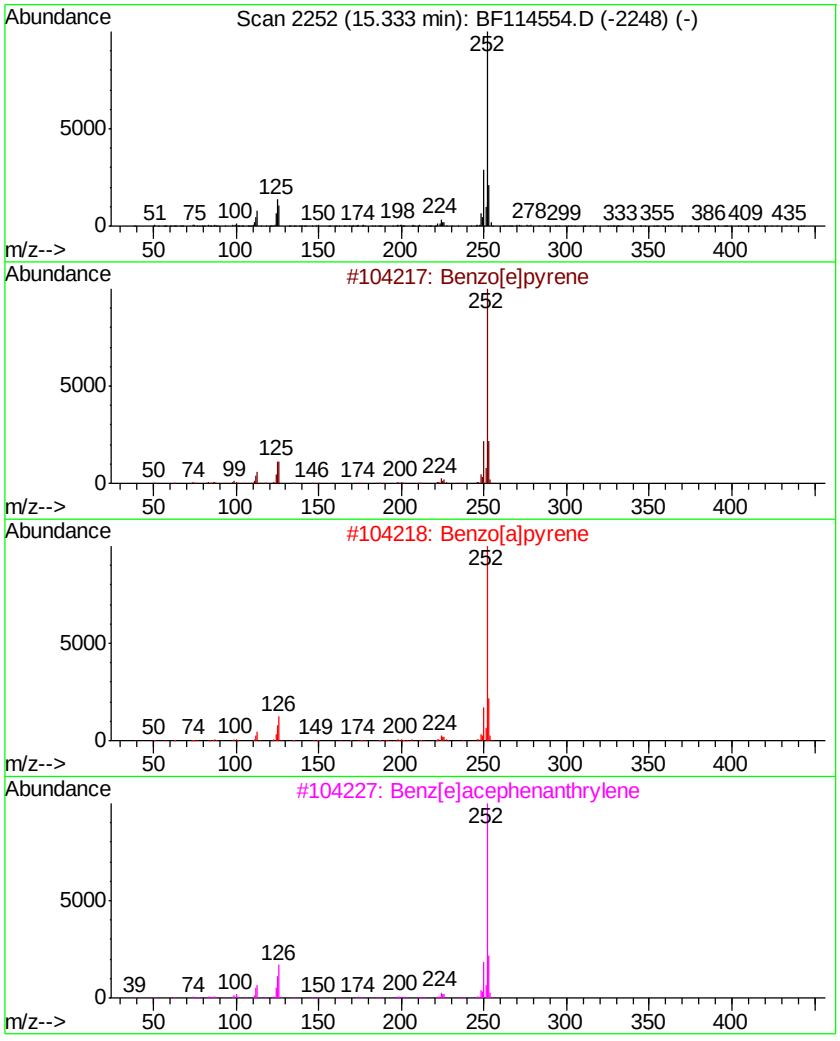
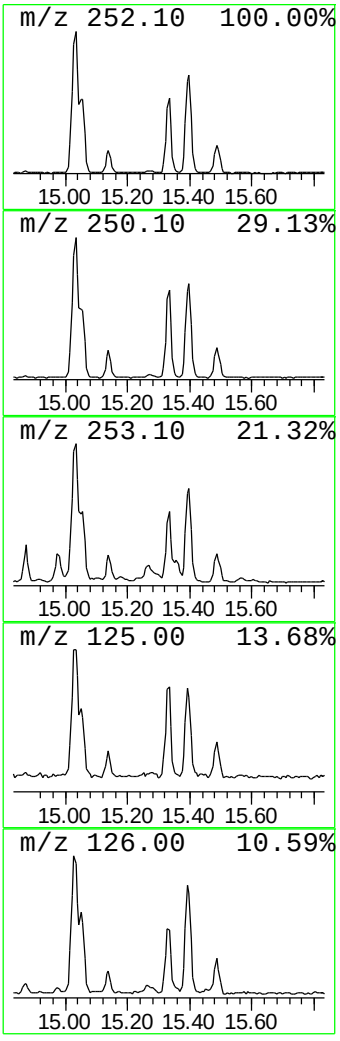
Peak Number 13 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.33	19.11 ng	679929	Perylene-d12	15.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2	Benzo[a]pyrene	252	C20H12	000050-32-8	97
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	90
4	Perylene	252	C20H12	000198-55-0	90
5	Benzo[k]fluoranthene	252	C20H12	000207-08-9	74





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF052319\
Data File : BF114554.D
Acq On : 24 May 2019 5:04
Operator : JU/SJ
Sample : K3008-01 2X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : COMP-19

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.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
unknown6.59	6.59	20.8	ng	1067330	1	6.81	1024350	20.0
Silane, chlorotri...	10.85	2.2	ng	173407	4	11.34	1551770	20.0
Acetamide, N-(3-m...	11.00	2.1	ng	161111	4	11.34	1551770	20.0
Phenanthrene, 2-m...	11.87	3.1	ng	242703	4	11.34	1551770	20.0
4H-Cyclopenta[def...	11.95	4.6	ng	354788	4	11.34	1551770	20.0
11H-Benzo[b]fluorene	13.11	4.1	ng	663690	5	13.99	3253660	20.0
Pyrene, 1-methyl-	13.22	3.5	ng	576136	5	13.99	3253660	20.0
Benzo[b]naphtho[2...	13.73	2.1	ng	336893	5	13.99	3253660	20.0
Cyclopenta[cd]pyrene	13.79	2.8	ng	460802	5	13.99	3253660	20.0
11H-Benzo[a]fluor...	13.84	2.2	ng	360209	5	13.99	3253660	20.0
Oxirane, 2,2'-[(1...	15.19	10.9	ng	387230	6	15.46	711581	20.0
Benzo[e]pyrene	15.33	19.1	ng	679929	6	15.46	711581	20.0