

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
 Data File : BF101779.D
 Acq On : 27 Dec 2017 22:49
 Operator : SJ/JU
 Sample : I7064-05 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SAMPLE-3

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.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	5.004	493	496	511	rBV	70636	130437	7.08%	0.536%
2	5.381	553	560	563	rBV4	8577	18734	1.02%	0.077%
3	5.416	563	566	598	rVV	956856	1764756	95.81%	7.246%
4	6.439	736	740	758	rBV	1028515	1807291	98.12%	7.420%
5	6.563	758	761	767	rBV	1644361	1698260	92.20%	6.973%
6	6.786	796	799	807	rBV	861909	866117	47.02%	3.556%
7	6.945	822	826	841	rBV	1273384	1383920	75.13%	5.682%
8	7.363	893	897	915	rBV	752884	1093988	59.39%	4.492%
9	8.069	1014	1017	1040	rVB	1023806	1152557	62.57%	4.732%
10	8.792	1137	1140	1146	rBV	35968	45126	2.45%	0.185%
11	8.886	1153	1156	1162	rBV	17663	21876	1.19%	0.090%
12	9.139	1196	1199	1209	rBV	1836392	1841943	100.00%	7.563%
13	9.210	1209	1211	1215	rVV	53477	46510	2.53%	0.191%
14	9.251	1215	1218	1223	rVV	20631	24549	1.33%	0.101%
15	9.422	1241	1247	1251	rBV2	17861	32087	1.74%	0.132%
16	9.486	1256	1258	1261	rVV3	21478	25488	1.38%	0.105%
17	9.545	1261	1268	1274	rVV	135871	181997	9.88%	0.747%
18	9.616	1277	1280	1285	rVB5	12222	24487	1.33%	0.101%
19	9.692	1285	1293	1298	rBV	39008	53177	2.89%	0.218%
20	9.739	1298	1301	1304	rVB	97640	70432	3.82%	0.289%
21	9.827	1312	1316	1320	rBV	1052549	1008516	54.75%	4.141%
22	9.857	1320	1321	1329	rVB	55725	58168	3.16%	0.239%
23	10.039	1347	1352	1354	rBV2	35222	48365	2.63%	0.199%
24	10.210	1378	1381	1383	rBV	24726	20916	1.14%	0.086%
25	10.239	1383	1386	1388	rVB	98474	79485	4.32%	0.326%
26	10.380	1403	1410	1416	rVB2	36901	48472	2.63%	0.199%
27	10.457	1416	1423	1426	rBV3	31727	47987	2.61%	0.197%
28	10.621	1448	1451	1463	rBV	730110	883012	47.94%	3.625%
29	10.710	1463	1466	1475	rVB	98929	121133	6.58%	0.497%
30	10.904	1497	1499	1502	rBV	35901	42964	2.33%	0.176%
31	11.163	1535	1543	1550	rBV3	79918	133356	7.24%	0.548%
32	11.221	1550	1553	1558	rVB2	28468	32228	1.75%	0.132%
33	11.321	1566	1570	1572	rBV	938088	928845	50.43%	3.814%
34	11.345	1572	1574	1579	rVV	402816	368502	20.01%	1.513%

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35	11.404	1580	1584	1591	rVV	104109	127274	6.91%	0.523%
36	11.468	1591	1595	1597	rVV2	17751	20208	1.10%	0.083%
37	11.568	1608	1612	1613	rBV2	40971	39035	2.12%	0.160%
38	11.827	1652	1656	1659	rBV2	137045	156483	8.50%	0.642%
39	11.851	1659	1660	1667	rVV2	83572	89222	4.84%	0.366%
40	11.904	1667	1669	1671	rVV	31176	27528	1.49%	0.113%
41	11.939	1671	1675	1677	rVV2	93622	109322	5.94%	0.449%
42	11.957	1677	1678	1682	rVV	43986	40431	2.20%	0.166%
43	11.992	1682	1684	1687	rVB	28566	21865	1.19%	0.090%
44	12.115	1702	1705	1710	rBV	46859	45982	2.50%	0.189%
45	12.174	1713	1715	1719	rVB5	27878	35474	1.93%	0.146%
46	12.374	1745	1749	1751	rBV2	81721	90782	4.93%	0.373%
47	12.480	1765	1767	1772	rVB2	52057	59618	3.24%	0.245%
48	12.539	1773	1777	1782	rBV	745685	685952	37.24%	2.816%
49	12.615	1788	1790	1792	rBV3	37070	31301	1.70%	0.129%
50	12.692	1800	1803	1809	rBV4	42917	59290	3.22%	0.243%
51	12.751	1810	1813	1814	rVV	157443	146773	7.97%	0.603%
52	12.768	1814	1816	1822	rVV	666420	607655	32.99%	2.495%
53	12.821	1822	1825	1829	rVB2	40359	52526	2.85%	0.216%
54	12.898	1834	1838	1843	rVV	1549508	1331121	72.27%	5.465%
55	12.945	1844	1846	1849	rVV	32081	34436	1.87%	0.141%
56	12.992	1850	1854	1859	rVB2	44093	79560	4.32%	0.327%
57	13.098	1866	1872	1877	rBV2	99373	157028	8.53%	0.645%
58	13.162	1881	1883	1886	rVV2	66735	60901	3.31%	0.250%
59	13.204	1888	1890	1897	rVB3	67803	92364	5.01%	0.379%
60	13.298	1904	1906	1909	rBV	44011	39226	2.13%	0.161%
61	13.327	1909	1911	1914	rBV	49495	51860	2.82%	0.213%
62	13.362	1915	1917	1921	rVB	74153	69805	3.79%	0.287%
63	13.780	1985	1988	1993	rBV3	233520	289240	15.70%	1.188%
64	13.915	2009	2011	2014	rVB	75791	60677	3.29%	0.249%
65	13.968	2015	2020	2023	rBV2	946007	1300171	70.59%	5.338%
66	13.998	2023	2025	2030	rVB	386056	382444	20.76%	1.570%
67	14.080	2037	2039	2044	rBV2	65930	67239	3.65%	0.276%
68	15.021	2195	2199	2202	rBV	386333	550642	29.89%	2.261%
69	15.321	2247	2250	2255	rBV	183629	228234	12.39%	0.937%
70	15.386	2258	2261	2266	rVB	271468	356000	19.33%	1.462%
71	15.445	2267	2271	2275	rBV	523248	652584	35.43%	2.679%

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

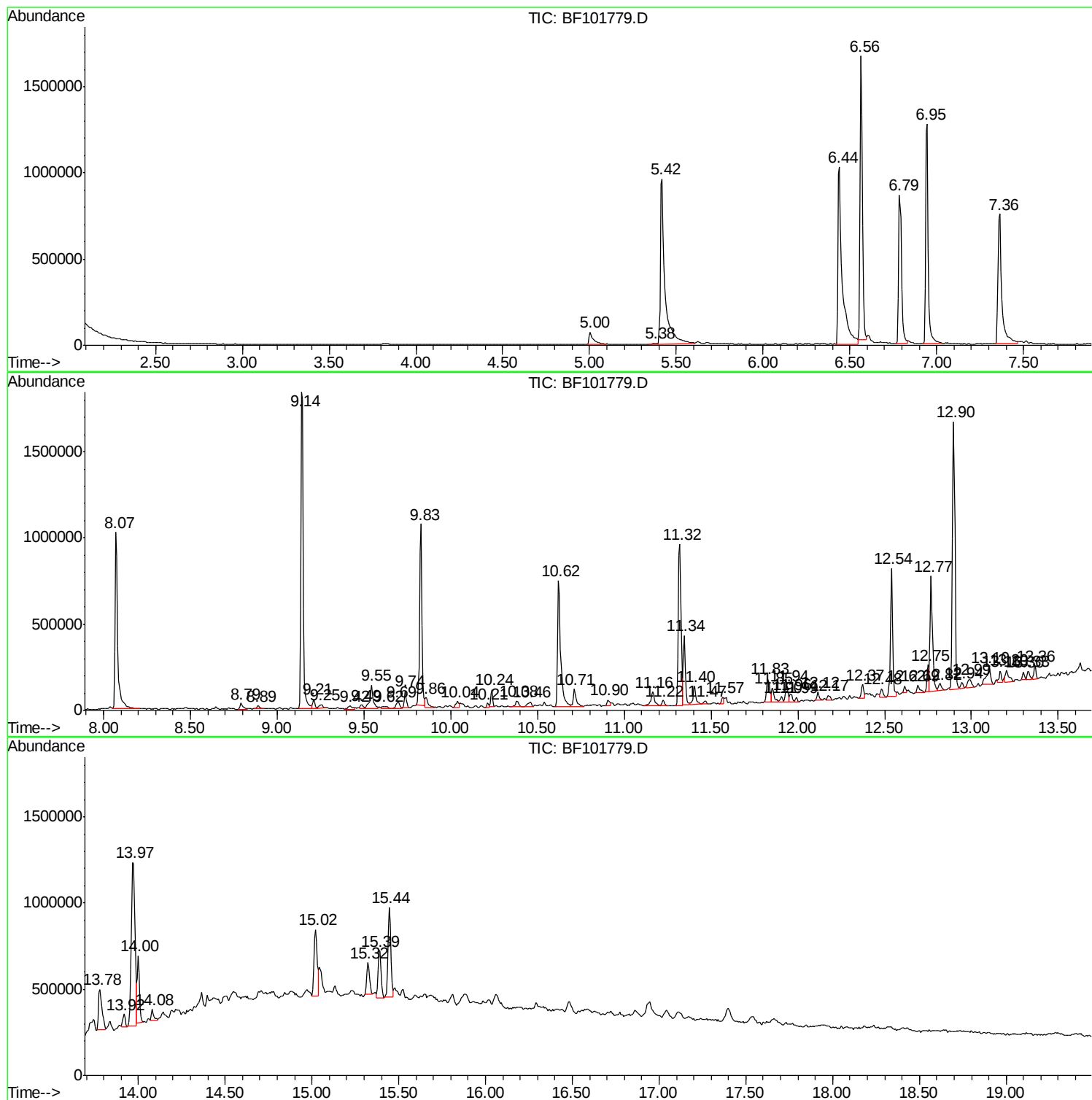
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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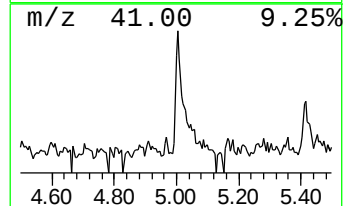
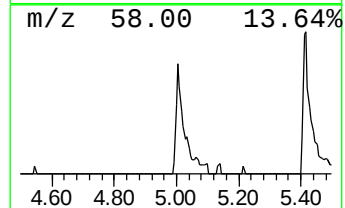
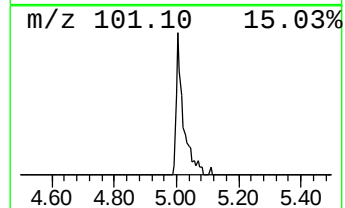
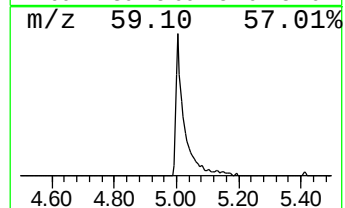
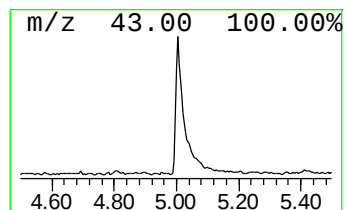
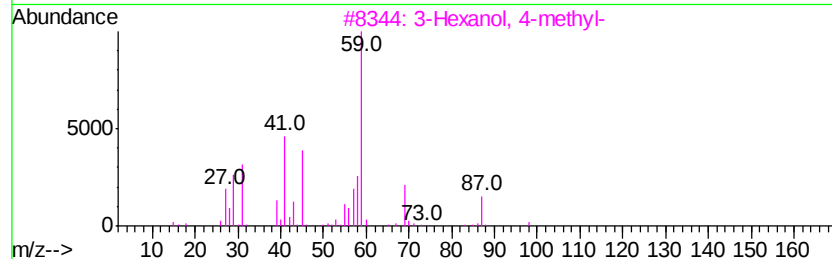
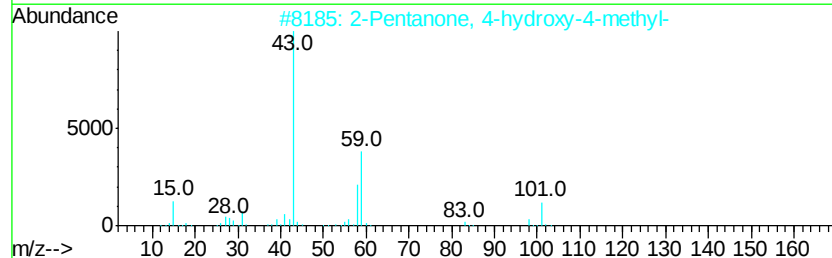
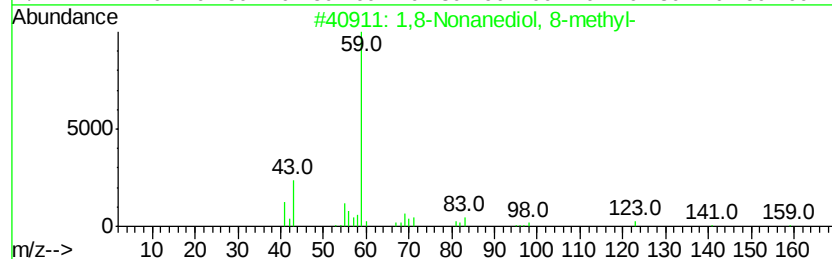
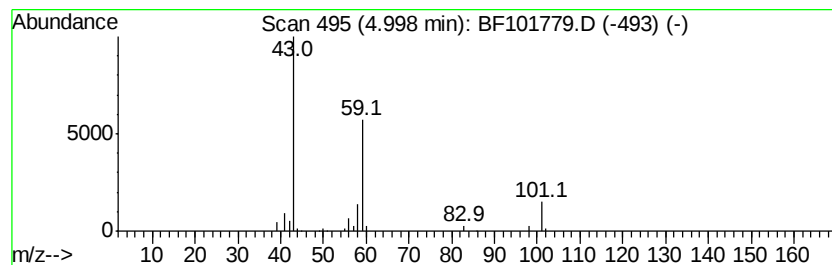
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown5.00 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	3.01 ng	130437	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	37
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	33
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4		3-Hexanol	102	C6H14O	000623-37-0	9
5		4-Pentene-2-ol, 2-methyl	100	C6H12O	000624-97-5	9



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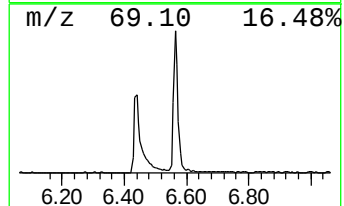
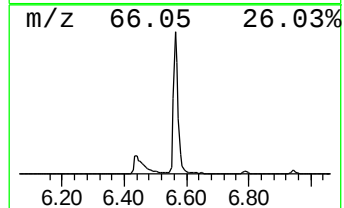
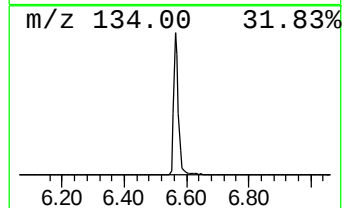
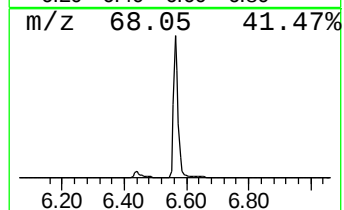
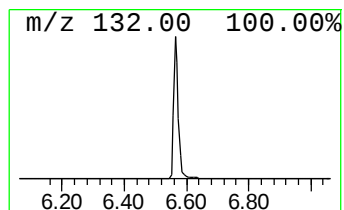
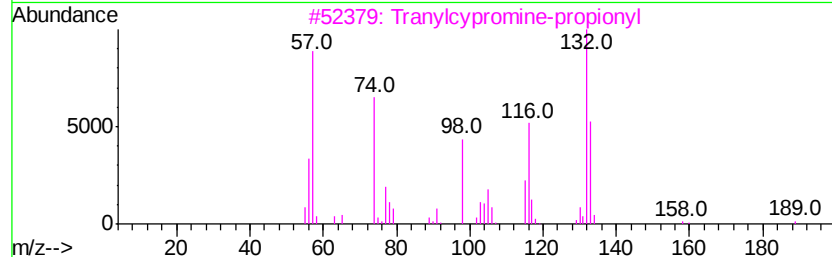
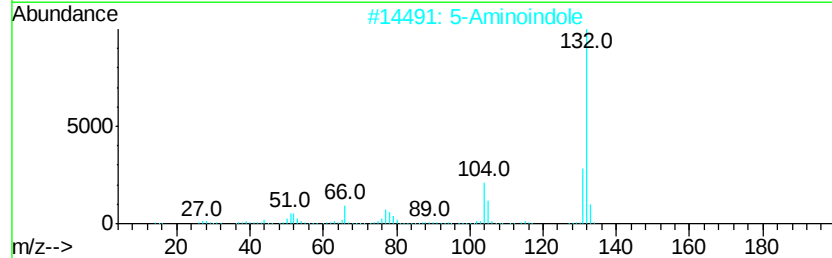
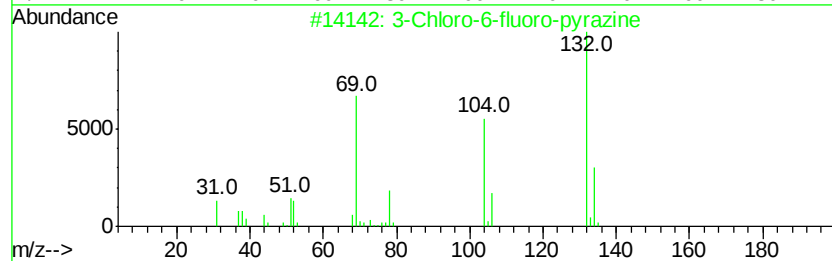
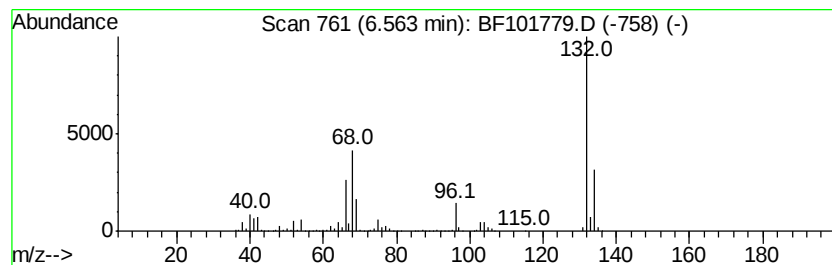
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	39.22 ng	1698260	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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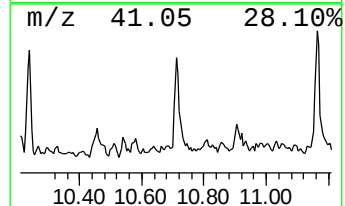
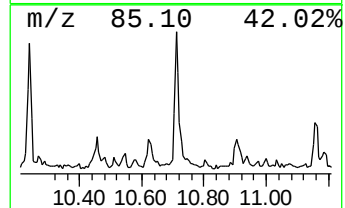
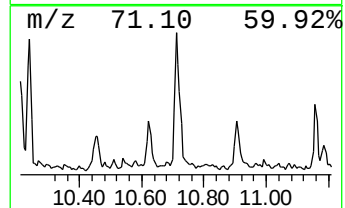
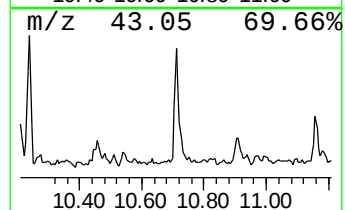
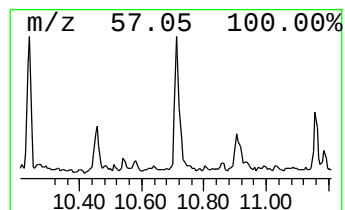
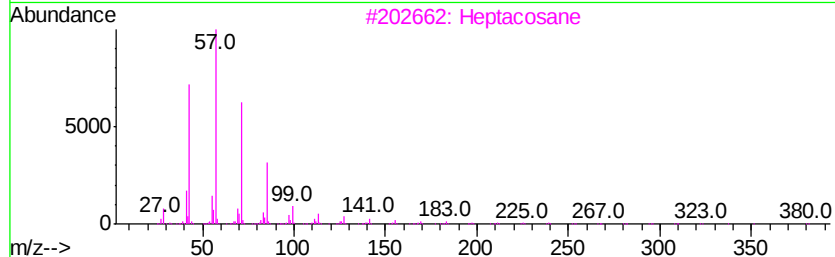
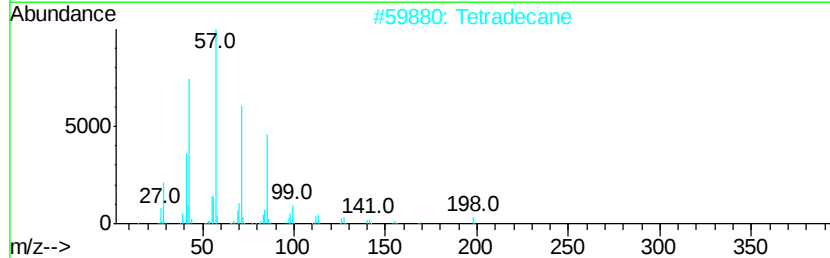
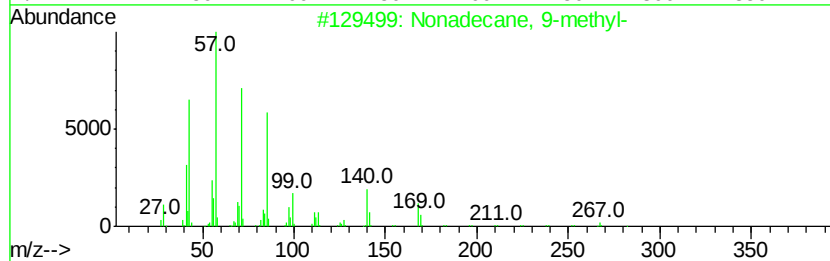
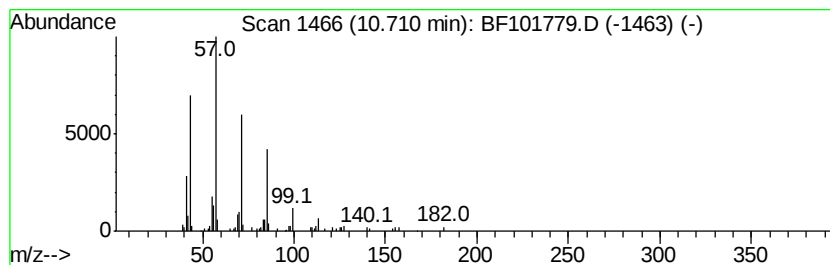
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Nonadecane, 9-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.71	2.61 ng	121133	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	86
2		Tetradecane	198	C14H30	000629-59-4	86
3		Heptacosane	380	C27H56	000593-49-7	86
4		Hexadecane	226	C16H34	000544-76-3	80
5		Pentadecane	212	C15H32	000629-62-9	72



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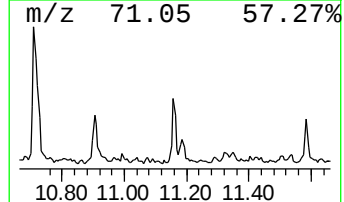
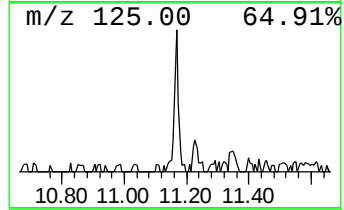
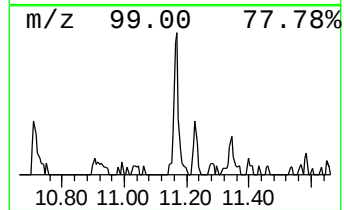
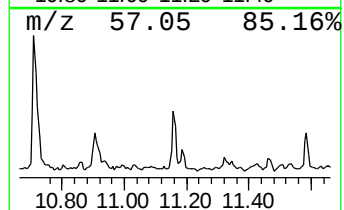
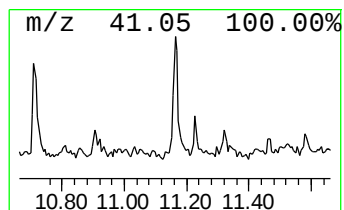
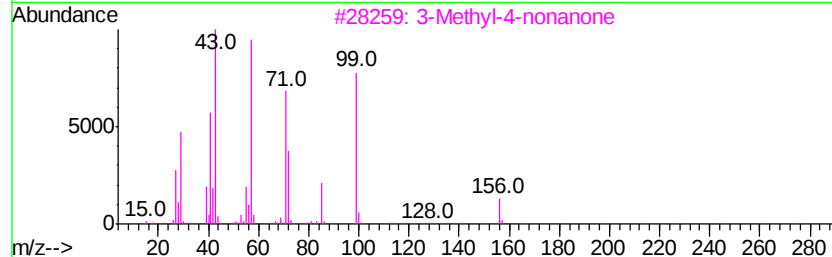
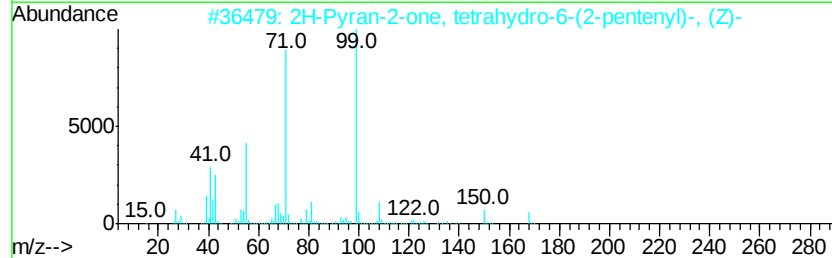
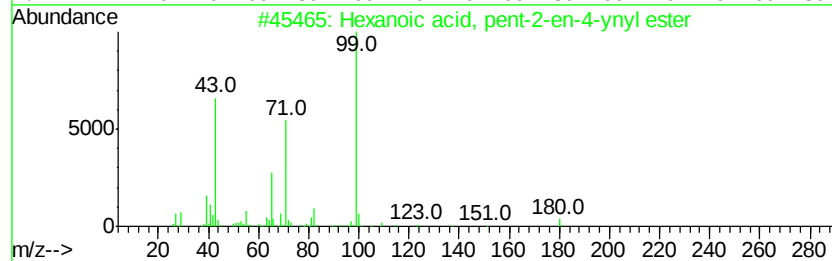
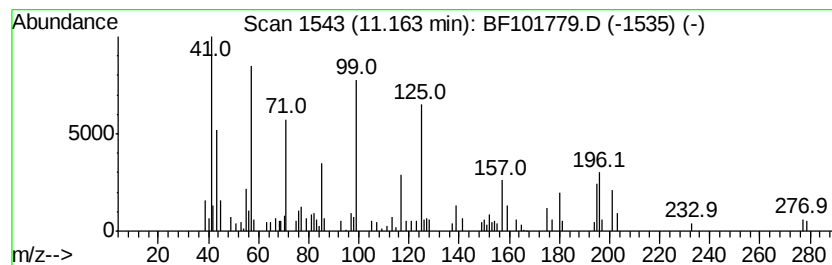
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown11.16 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.16	2.87 ng	133356	Phenanthrene-d10	11.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexanoic acid, pent-2-en-4-ynyl ...	180	C11H16O2	1000299-35-3	35
2		2H-Pyran-2-one, tetrahydro-6-(2-...	168	C10H16O2	025524-95-2	27
3		3-Methyl-4-nonanone	156	C10H20O	035778-39-3	27
4		Hexanoic acid, 2-propenyl ester	156	C9H16O2	000123-68-2	27
5		Hexanoic acid, 2-bromo-4-fluorop...	288	C12H14BrFO2	003259-26-5	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122717\
Data File : BF101779.D
Acq On : 27 Dec 2017 22:49
Operator : SJ/JU
Sample : I7064-05 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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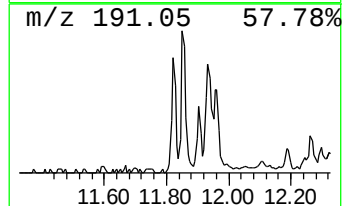
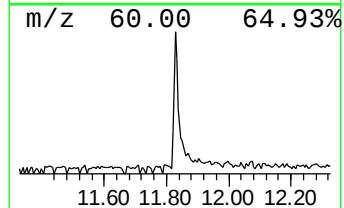
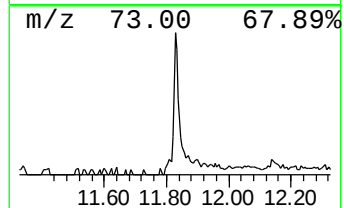
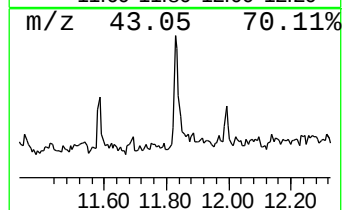
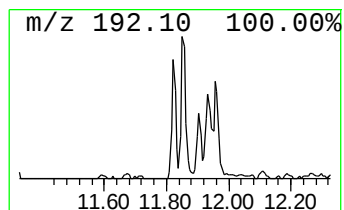
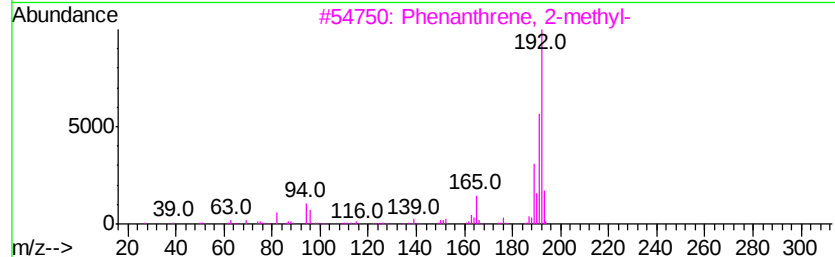
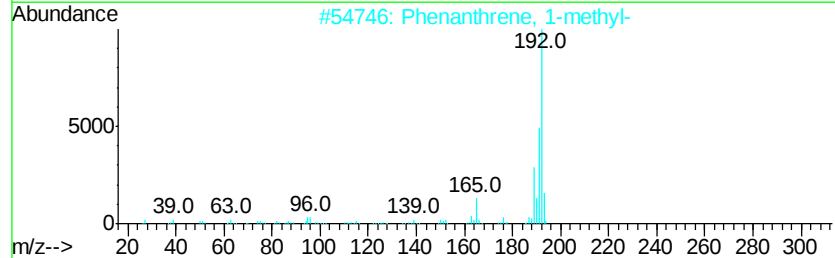
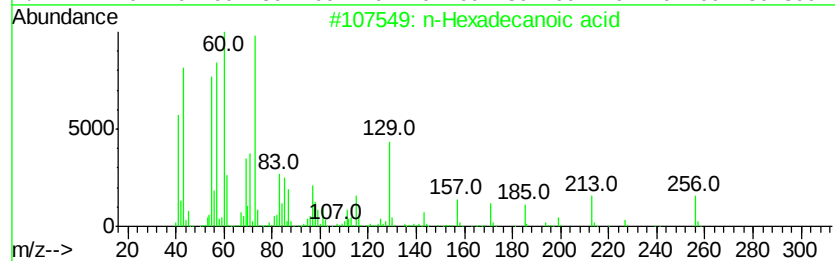
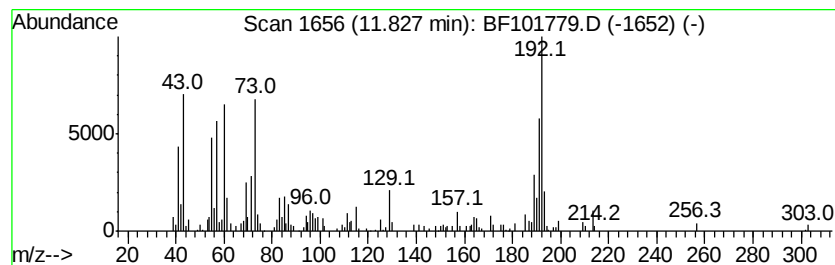
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.83	3.37 ng	156483	Phenanthrene-d10	11.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	78
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	78
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	70
5		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	70



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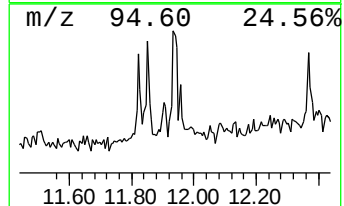
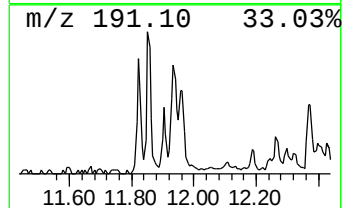
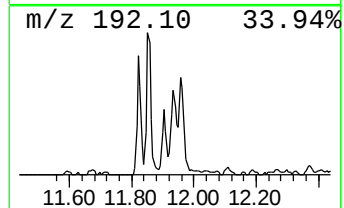
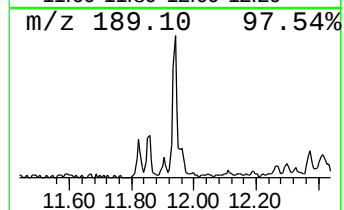
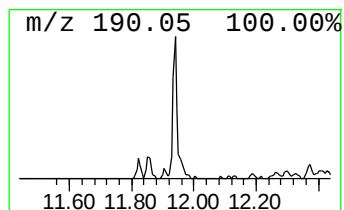
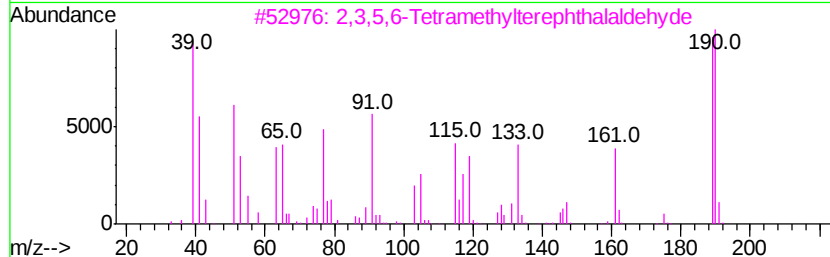
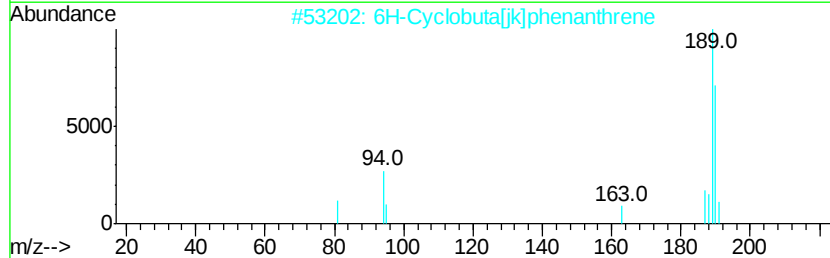
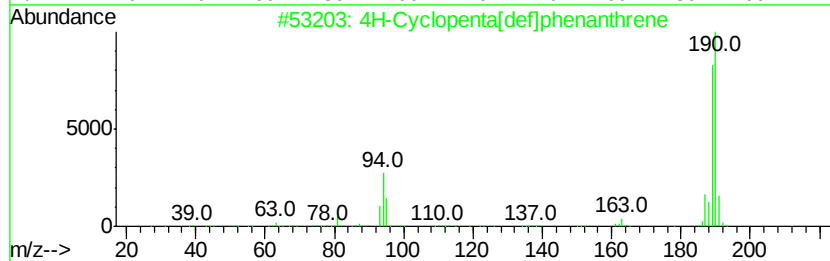
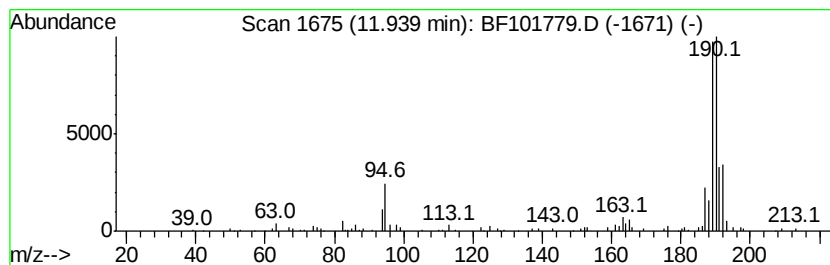
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.94	2.35 ng	109322	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	72
3		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50
4		3-Bromo-2-furoic acid	190	C5H3BrO3	014903-90-3	10
5		Propanoic acid, 2-(2,4-dichlorop...	378	C15H10Cl4O3	284488-02-4	10



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Misc :
ALS Vial : 20 Sample Multiplier: 1

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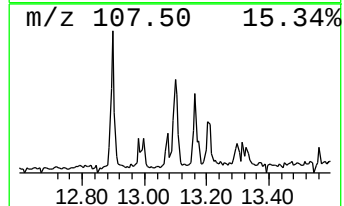
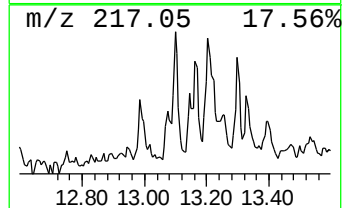
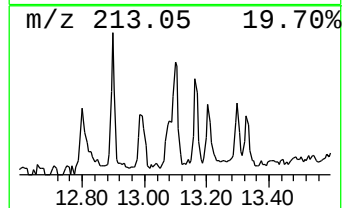
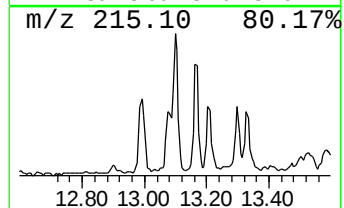
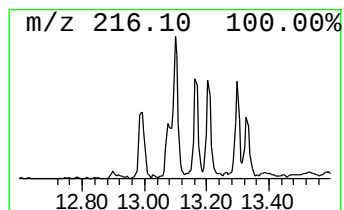
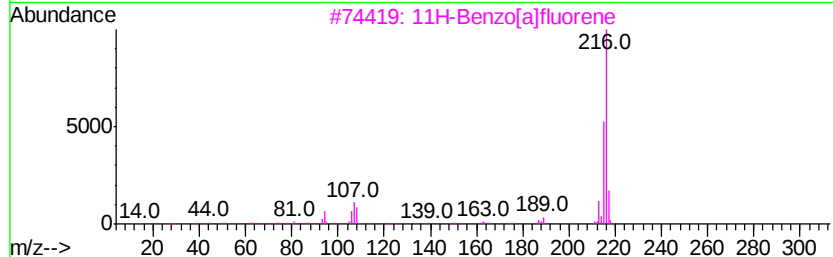
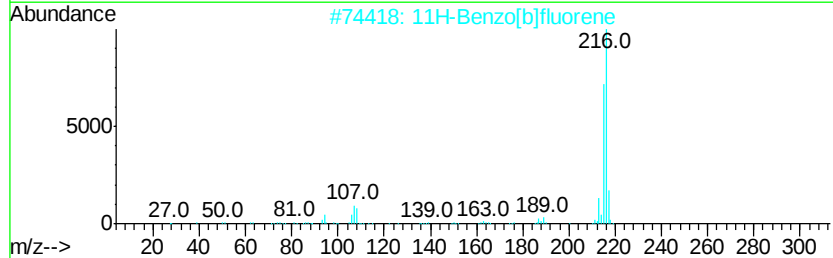
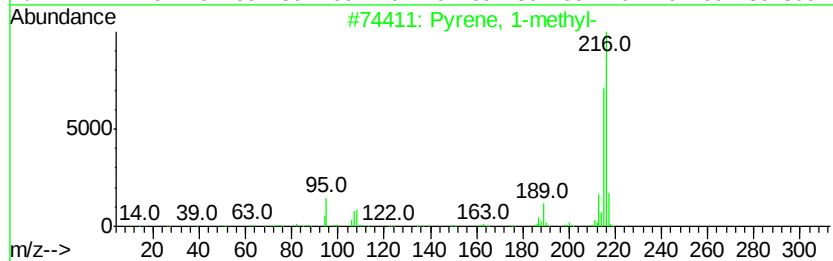
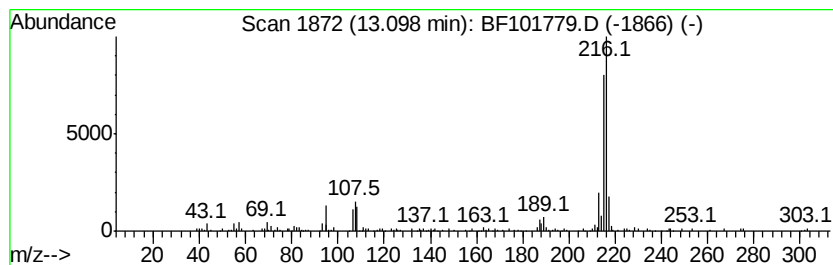
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pyrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.10	2.42 ng	157028	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
2		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
5		Pyrene, 2-methyl-	216	C17H12	003442-78-2	89



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122717\
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Operator : SJ/JU
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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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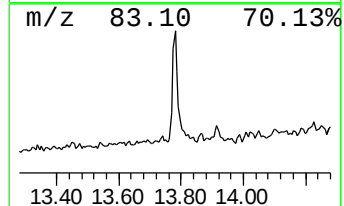
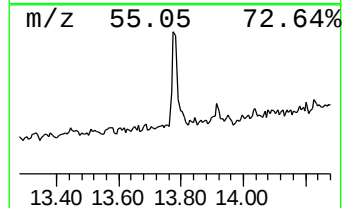
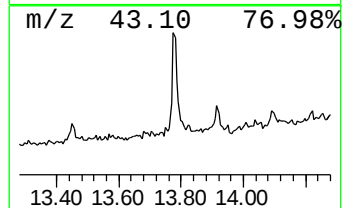
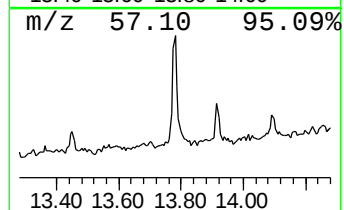
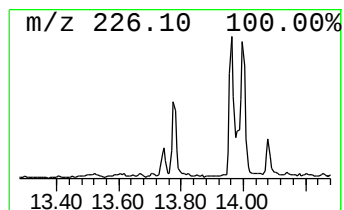
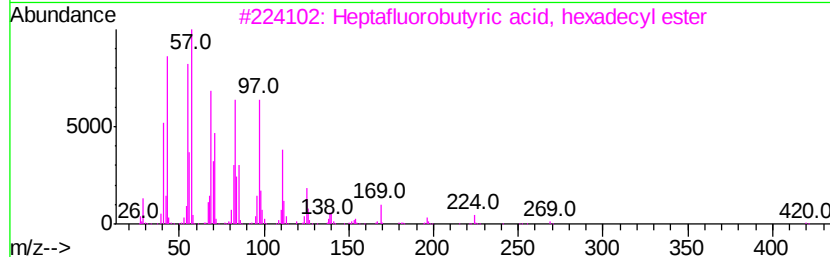
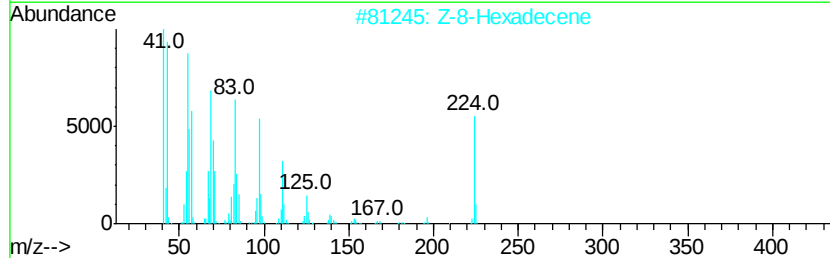
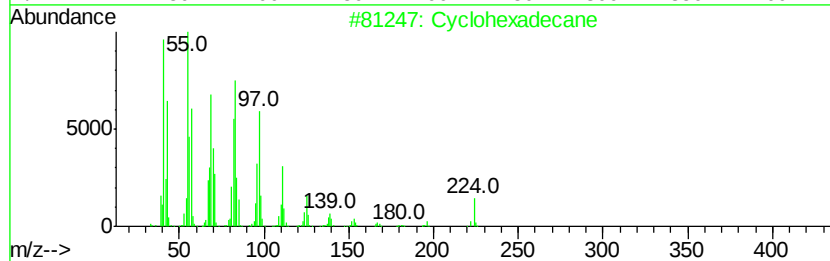
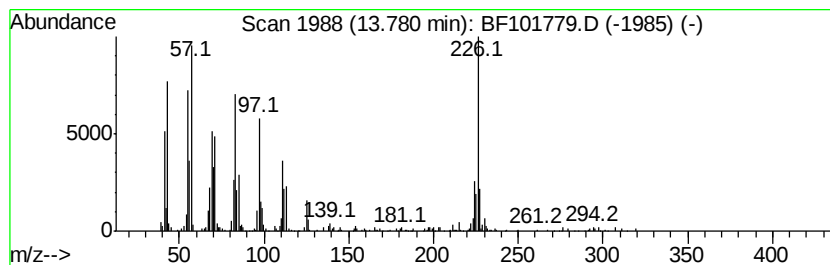
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Cyclohexadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.78	4.45 ng	289240	Chrysene-d12	13.97

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexadecane	224	C16H32	000295-65-8	51
2		Z-8-Hexadecene	224	C16H32	1000130-87-5	42
3		Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	42
4		1-Octadecanol	270	C18H38O	000112-92-5	38
5		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
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Acq On : 27 Dec 2017 22:49
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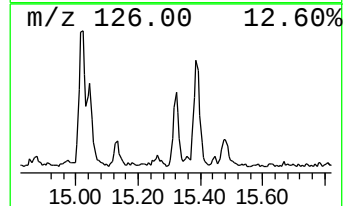
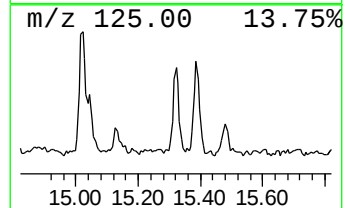
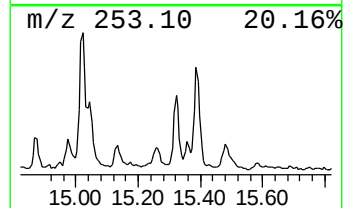
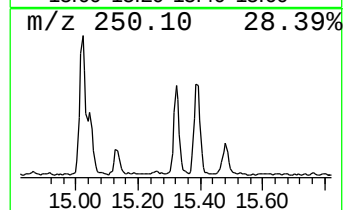
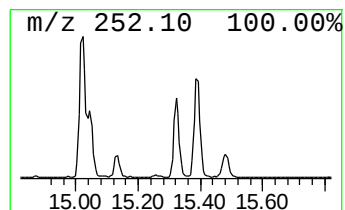
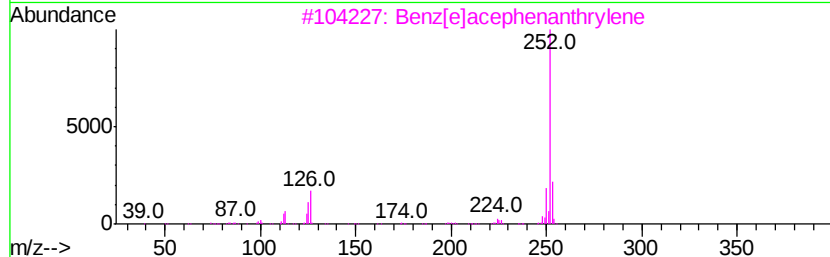
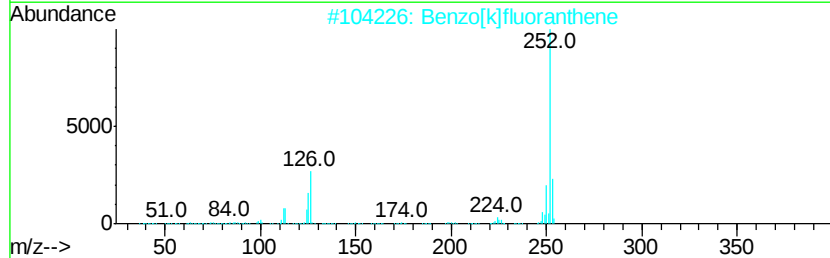
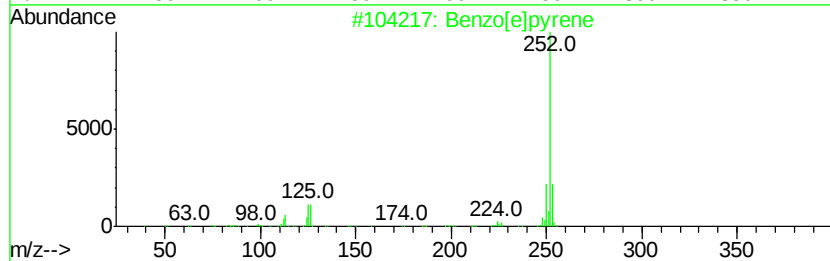
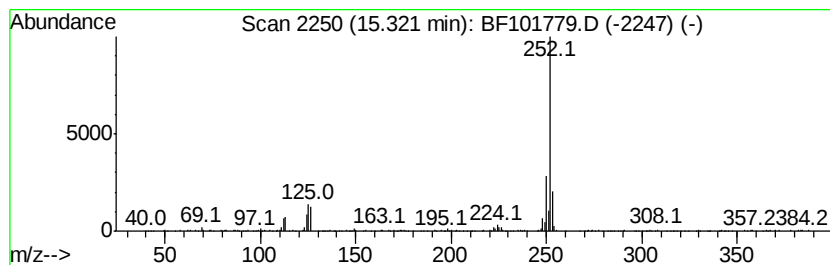
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.32	6.99 ng	228234	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	96
4		Benzo[i]fluoranthene	252	C20H12	000205-82-3	95
5		Benzo[a]pyrene	252	C20H12	000050-32-8	95



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Instrument :
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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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unknown6.56	6.56	39.2	ng	1698260	1	6.79	866117	20.0
Nonadecane, 9-met...	10.71	2.6	ng	121133	4	11.32	928845	20.0
unknown11.16	11.16	2.9	ng	133356	4	11.32	928845	20.0
n-Hexadecanoic acid	11.83	3.4	ng	156483	4	11.32	928845	20.0
4H-Cyclopenta[def...	11.94	2.4	ng	109322	4	11.32	928845	20.0
Pyrene, 1-methyl-	13.10	2.4	ng	157028	5	13.97	1300170	20.0
Cyclohexadecane	13.78	4.5	ng	289240	5	13.97	1300170	20.0
Benzo[e]pyrene	15.32	7.0	ng	228234	6	15.44	652584	20.0