

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061218\
 Data File : BF106344.D
 Acq On : 12 Jun 2018 15:07
 Operator : JU/SJ
 Sample : J3245-03 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-061118-B

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061118.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	5.195	482	486	492	rVB	155383	150500	7.90%	0.539%
2	5.589	549	553	556	rBV	1763388	1504735	78.96%	5.393%
3	6.589	718	723	729	rBV	1645229	1489106	78.14%	5.337%
4	6.736	744	748	752	rVB	1782133	1523613	79.95%	5.460%
5	6.972	784	788	791	rVB	1191991	922612	48.41%	3.306%
6	7.125	811	814	818	rVB	1451062	1213820	63.69%	4.350%
7	7.531	879	883	886	rBV	1062120	912917	47.90%	3.272%
8	8.254	1002	1006	1009	rBV	1352963	1078187	56.58%	3.864%
9	8.966	1123	1127	1131	rVB	76829	57326	3.01%	0.205%
10	9.066	1141	1144	1147	rBV	59208	44638	2.34%	0.160%
11	9.325	1184	1188	1191	rBV	1958544	1516671	79.58%	5.435%
12	9.589	1230	1233	1238	rBV	32631	43121	2.26%	0.155%
13	9.666	1242	1246	1249	rBV	58305	56365	2.96%	0.202%
14	9.719	1252	1255	1261	rVB	226991	195189	10.24%	0.700%
15	9.783	1261	1266	1267	rBV	31537	24226	1.27%	0.087%
16	9.872	1277	1281	1285	rBV	108064	91084	4.78%	0.326%
17	9.924	1286	1290	1294	rVB3	37011	40362	2.12%	0.145%
18	10.013	1301	1305	1308	rVV	1161672	955208	50.12%	3.423%
19	10.042	1308	1310	1313	rVB	134705	102186	5.36%	0.366%
20	10.107	1318	1321	1326	rBV4	13708	22240	1.17%	0.080%
21	10.213	1333	1339	1342	rVB2	99499	99808	5.24%	0.358%
22	10.248	1342	1345	1349	rBV	33366	28609	1.50%	0.103%
23	10.324	1354	1358	1360	rBV2	33452	37169	1.95%	0.133%
24	10.354	1360	1363	1369	rVV	21428	25760	1.35%	0.092%
25	10.430	1369	1376	1379	rVB2	54232	85856	4.51%	0.308%
26	10.560	1394	1398	1403	rVB	166801	153594	8.06%	0.550%
27	10.624	1403	1409	1410	rBV4	20529	27957	1.47%	0.100%
28	10.648	1410	1413	1415	rVB	26162	27246	1.43%	0.098%
29	10.713	1422	1424	1428	rVB	24013	22365	1.17%	0.080%
30	10.795	1435	1438	1442	rBV	908419	812016	42.61%	2.910%
31	10.901	1453	1456	1457	rBV	49742	40344	2.12%	0.145%
32	10.919	1457	1459	1466	rVB2	48160	56831	2.98%	0.204%
33	11.101	1484	1490	1493	rBV4	53093	76913	4.04%	0.276%
34	11.136	1493	1496	1499	rVB2	23063	28364	1.49%	0.102%

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35	11.260	1514	1517	1520	rVB4	18424	21400	1.12%	0.077%
36	11.301	1522	1524	1527	rVV3	25819	25311	1.33%	0.091%
37	11.348	1529	1532	1534	rVV2	43824	39517	2.07%	0.142%
38	11.395	1538	1540	1544	rVB	74576	71176	3.73%	0.255%
39	11.501	1554	1558	1560	rBV	1174785	1006920	52.84%	3.609%
40	11.524	1560	1562	1566	rVV	999222	769406	40.37%	2.757%
41	11.577	1568	1571	1573	rVV	266003	213572	11.21%	0.765%
42	11.607	1573	1576	1579	rVB	38216	38612	2.03%	0.138%
43	11.730	1590	1597	1600	rBV	106240	112525	5.90%	0.403%
44	11.830	1611	1614	1620	rVB	52974	51756	2.72%	0.185%
45	11.913	1626	1628	1633	rVB	33457	34535	1.81%	0.124%
46	12.001	1640	1643	1646	rBV	143415	167958	8.81%	0.602%
47	12.030	1646	1648	1651	rVB	190890	147698	7.75%	0.529%
48	12.077	1653	1656	1659	rBV	66020	55994	2.94%	0.201%
49	12.118	1659	1663	1665	rBV2	241950	277658	14.57%	0.995%
50	12.183	1671	1674	1676	rBV3	30629	33837	1.78%	0.121%
51	12.236	1680	1683	1685	rVB2	29331	28270	1.48%	0.101%
52	12.295	1690	1693	1695	rVV	95025	83543	4.38%	0.299%
53	12.318	1695	1697	1701	rVB3	51192	59437	3.12%	0.213%
54	12.430	1713	1716	1717	rBV3	28332	25609	1.34%	0.092%
55	12.448	1717	1719	1721	rVV	51778	37172	1.95%	0.133%
56	12.477	1721	1724	1726	rVV	59717	50058	2.63%	0.179%
57	12.501	1726	1728	1730	rVB2	46388	34210	1.80%	0.123%
58	12.554	1734	1737	1739	rBV	123752	117895	6.19%	0.423%
59	12.589	1739	1743	1745	rVV3	60243	82026	4.30%	0.294%
60	12.636	1748	1751	1755	rBV2	59102	79934	4.19%	0.286%
61	12.718	1761	1765	1768	rBV	1310288	1116357	58.58%	4.001%
62	12.754	1768	1771	1773	rVV2	157392	135562	7.11%	0.486%
63	12.777	1773	1775	1777	rVV2	45683	38797	2.04%	0.139%
64	12.813	1778	1781	1783	rVB2	46803	44833	2.35%	0.161%
65	12.871	1788	1791	1793	rBV	71980	63016	3.31%	0.226%
66	12.948	1801	1804	1810	rVB	1143305	1041110	54.63%	3.731%
67	12.995	1810	1812	1816	rVB2	73104	74778	3.92%	0.268%
68	13.060	1821	1823	1824	rVV2	60946	48731	2.56%	0.175%
69	13.083	1824	1827	1834	rVB	2038685	1807754	94.86%	6.479%
70	13.160	1836	1840	1844	rBV3	96136	167018	8.76%	0.599%
71	13.248	1852	1855	1856	rBV3	63698	67313	3.53%	0.241%

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72	13.271	1856	1859	1862	rVB	240947	225022	11.81%	0.806%
73	13.301	1862	1864	1866	rBV3	35461	28109	1.47%	0.101%
74	13.342	1868	1871	1873	rBV	186993	163650	8.59%	0.586%
75	13.377	1874	1877	1880	rBV2	134145	131185	6.88%	0.470%
76	13.471	1891	1893	1896	rBV	103860	77626	4.07%	0.278%
77	13.501	1896	1898	1901	rVV	88970	83227	4.37%	0.298%
78	13.901	1964	1966	1968	rVB	107937	85823	4.50%	0.308%
79	13.930	1968	1971	1973	rVB2	72869	56703	2.98%	0.203%
80	13.971	1973	1978	1987	rBV5	230141	437862	22.98%	1.569%
81	14.160	2005	2010	2012	rBV2	1452556	1905725	100.00%	6.830%
82	14.183	2012	2014	2021	rVB	612111	555084	29.13%	1.989%
83	14.265	2026	2028	2030	rBV	88999	73576	3.86%	0.264%
84	15.248	2192	2195	2199	rBV	365525	572897	30.06%	2.053%
85	15.571	2248	2250	2256	rVB	247638	323286	16.96%	1.159%
86	15.642	2258	2262	2267	rBV	307789	409422	21.48%	1.467%
87	15.712	2269	2274	2278	rBV	709142	1035783	54.35%	3.712%

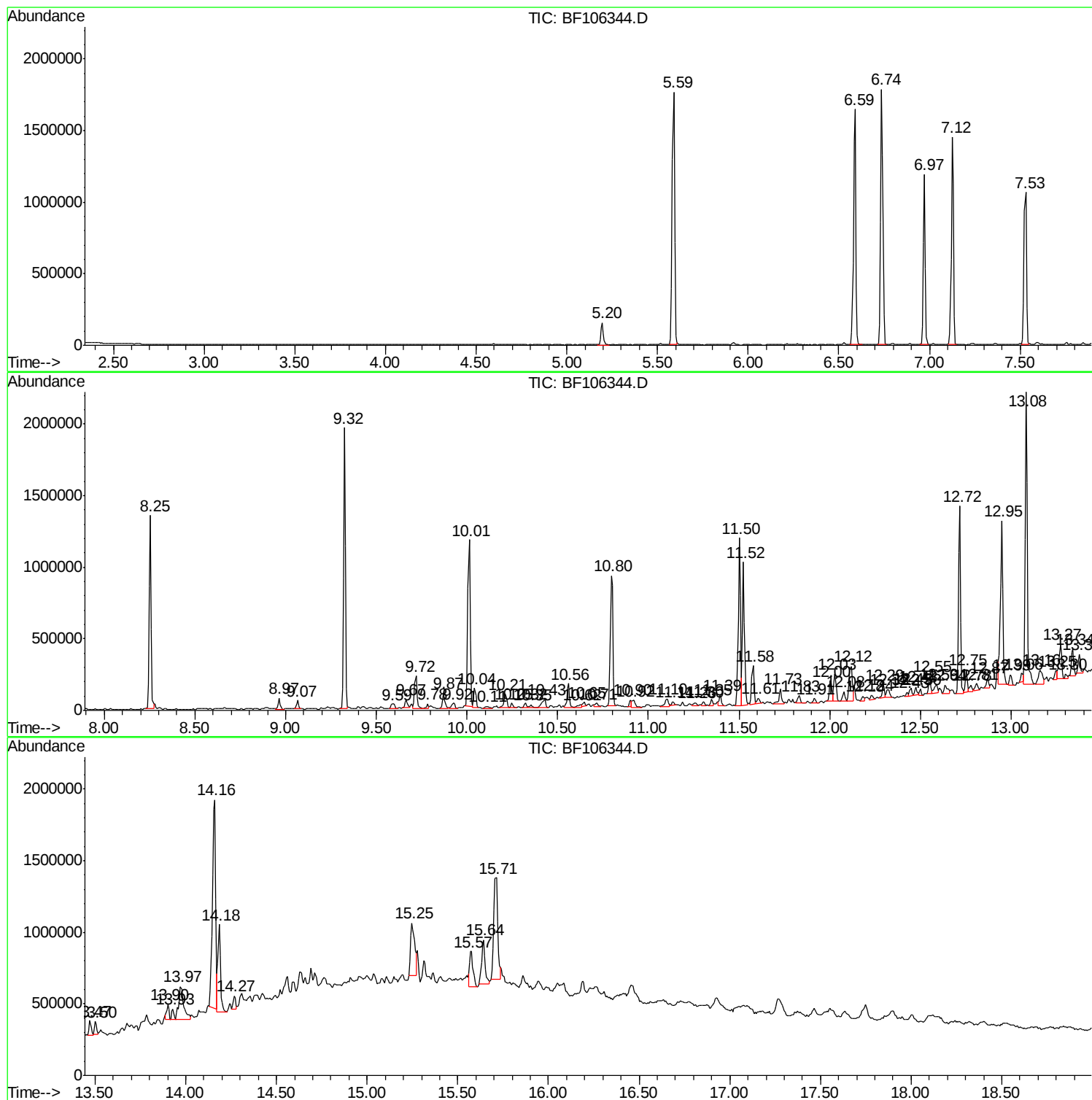
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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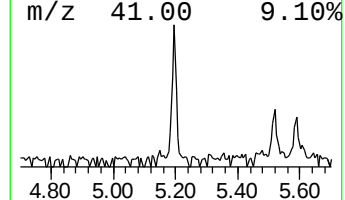
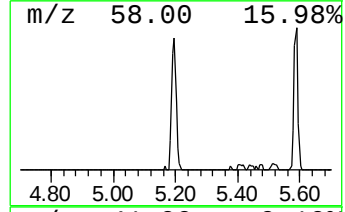
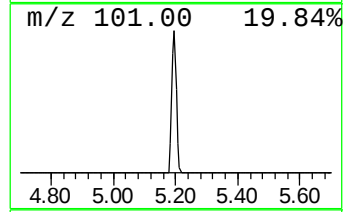
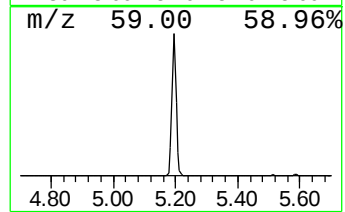
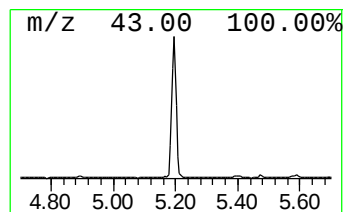
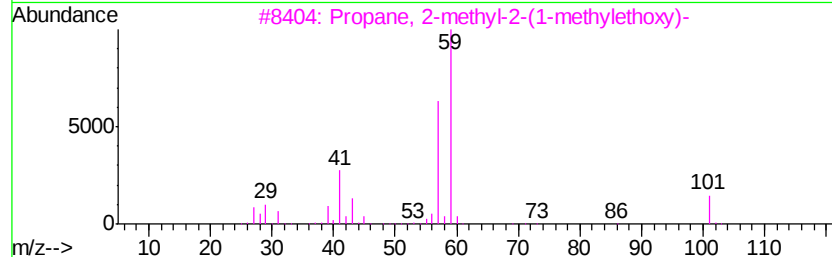
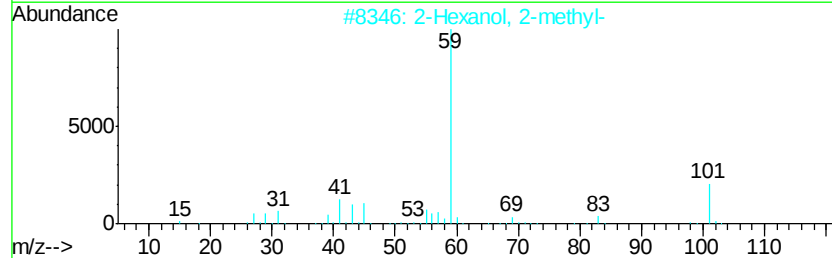
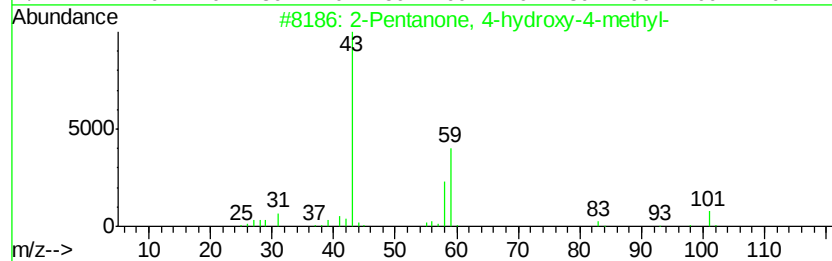
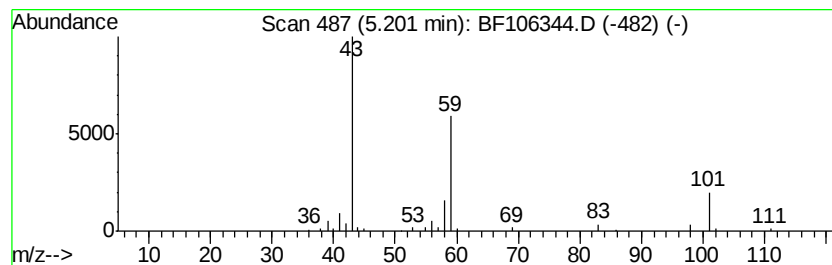
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	3.26 ng	150500	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
3			Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	36
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5			3-Octanol	130	C8H18O	000589-98-0	33



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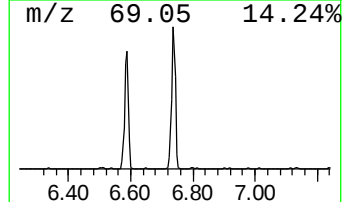
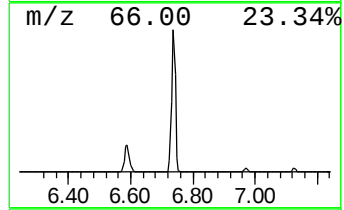
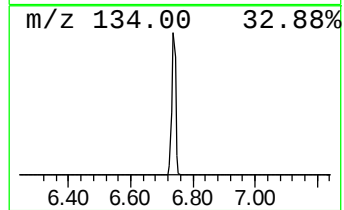
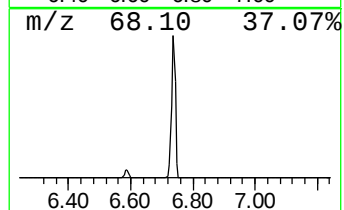
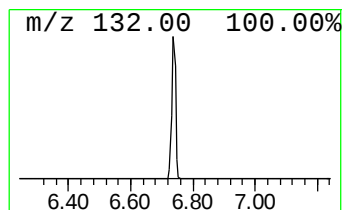
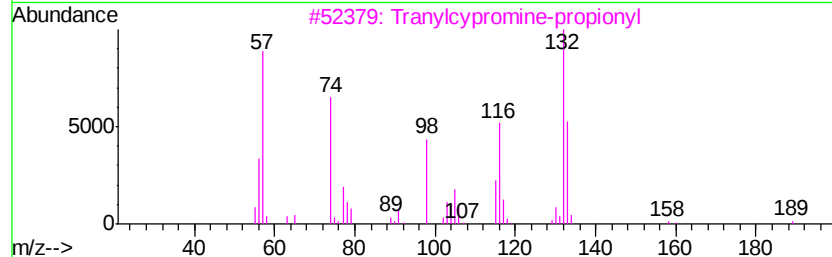
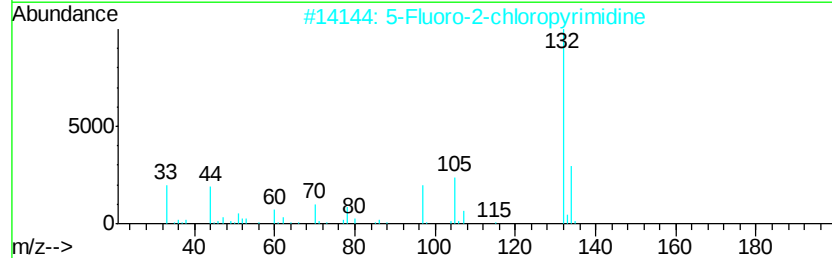
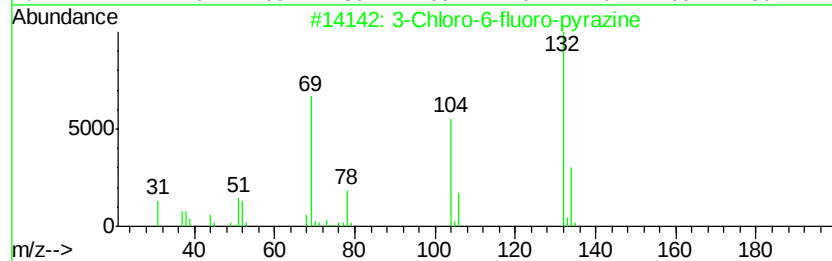
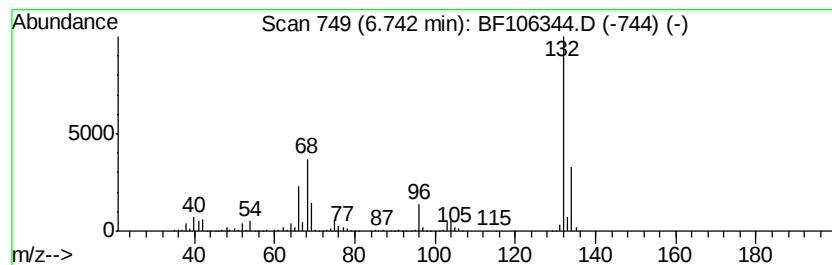
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	33.03 ng	1523610	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranvlcypramine-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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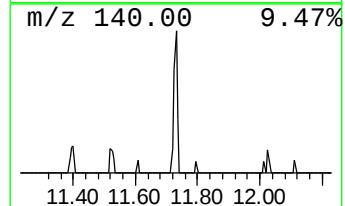
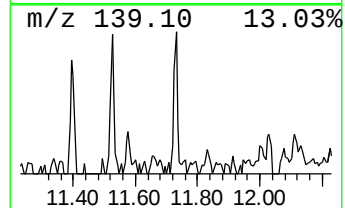
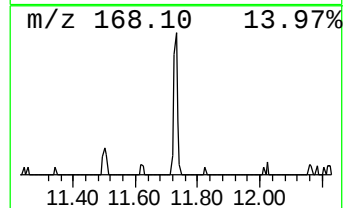
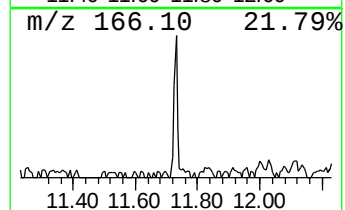
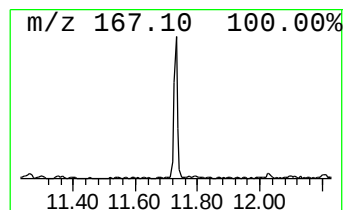
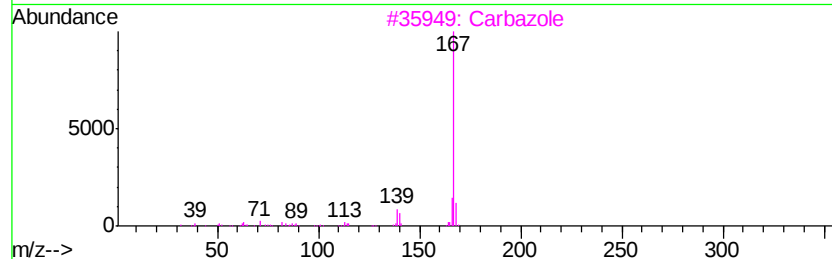
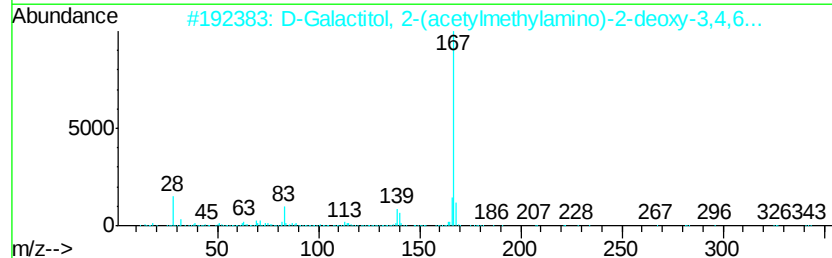
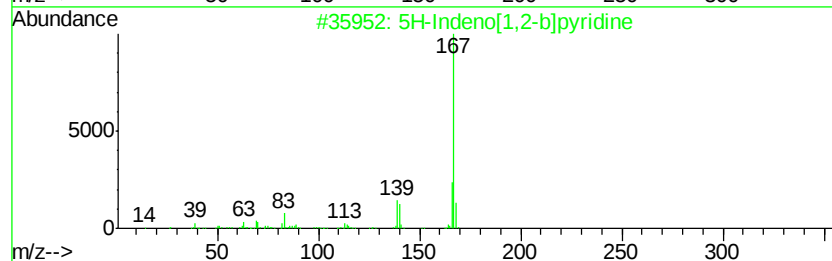
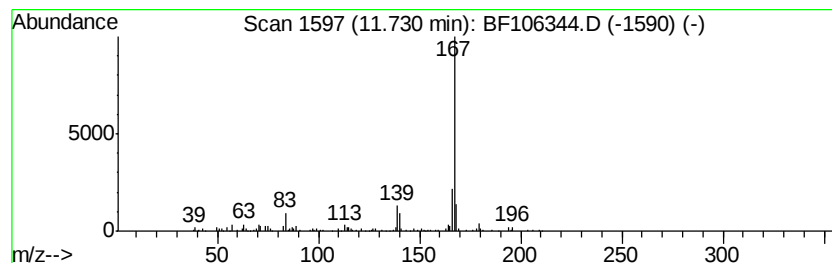
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 5H-Indeno[1,2-b]pyridine Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.73	2.24 ng	112525	Phenanthrene-d10	11.50

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	94
2		D-Galactitol, 2-(acetylmethylami...	363	C16H29N08	074410-42-7	87
3		Carbazole	167	C12H9N	000086-74-8	87
4		9H-Carbazole, 9-nitroso-	196	C12H8N2O	002788-23-0	80
5		9H-Carbazole-9-methanol	197	C13H11NO	002409-36-1	72



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CL-01-061118-B

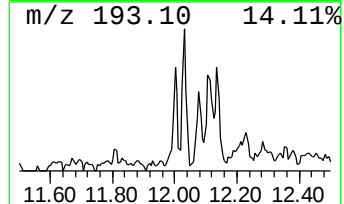
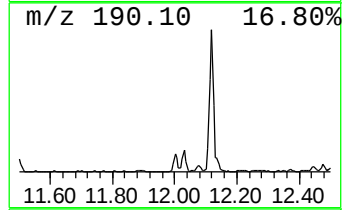
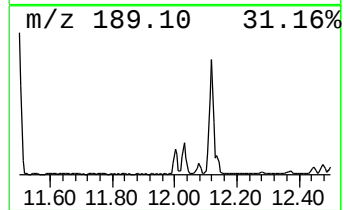
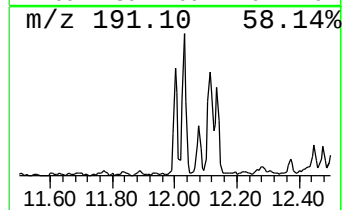
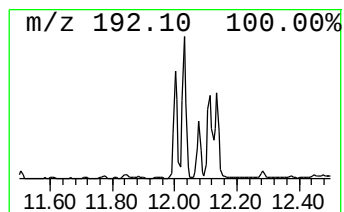
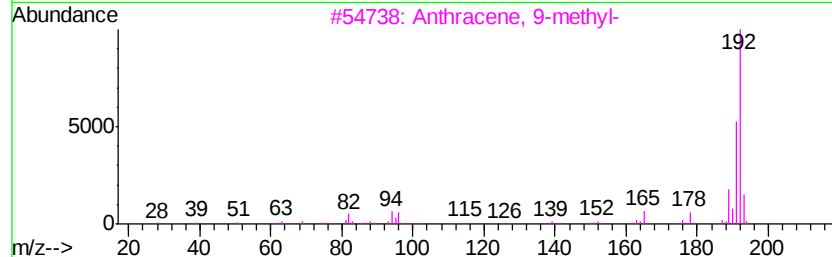
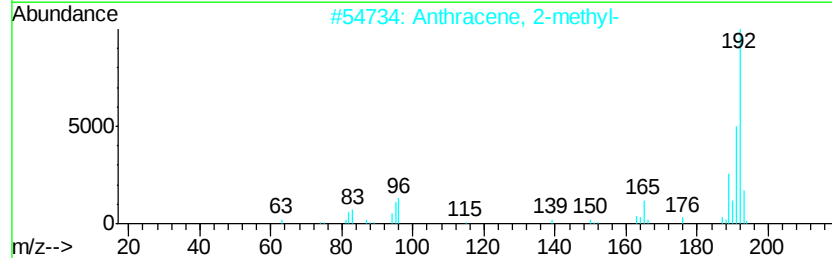
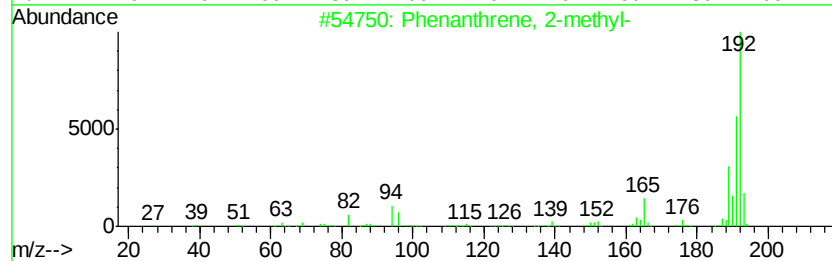
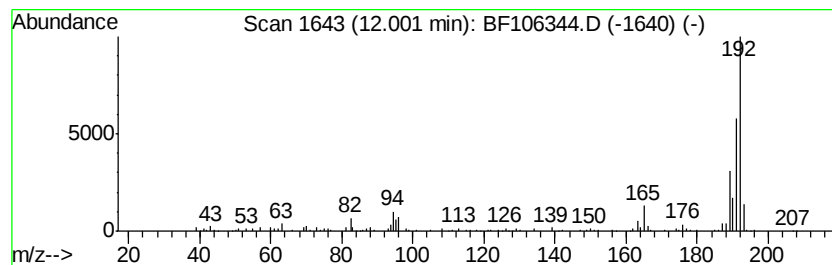
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	3.34 ng	167958	Phenanthrene-d10	11.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	87
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	87
4			1H-Cyclopropa[1,1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	87
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061218\
Data File : BF106344.D
Acq On : 12 Jun 2018 15:07
Operator : JU/SJ
Sample : J3245-03 2X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CL-01-061118-B

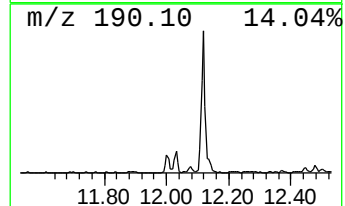
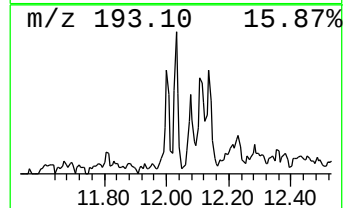
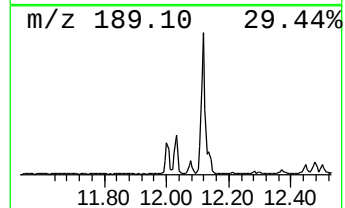
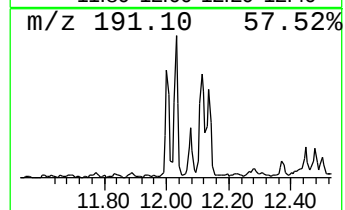
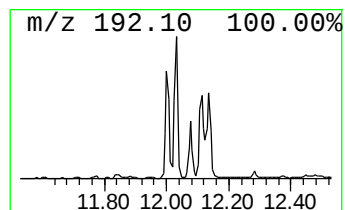
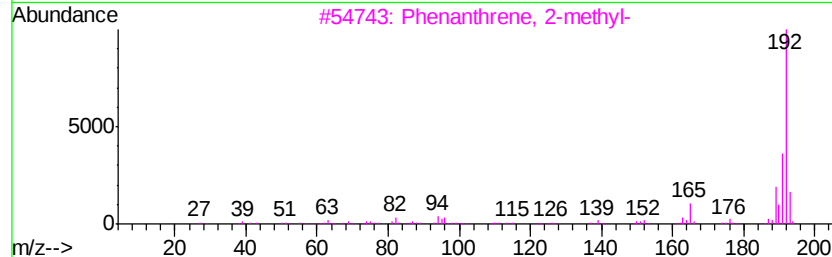
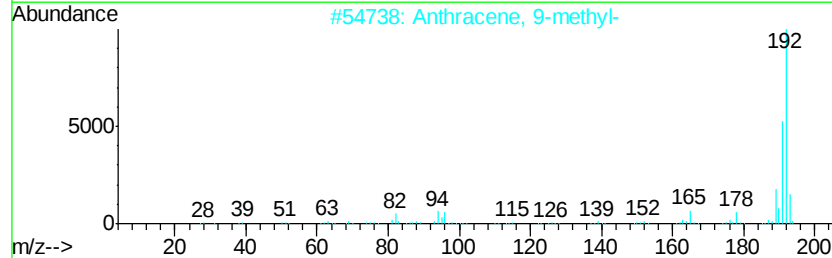
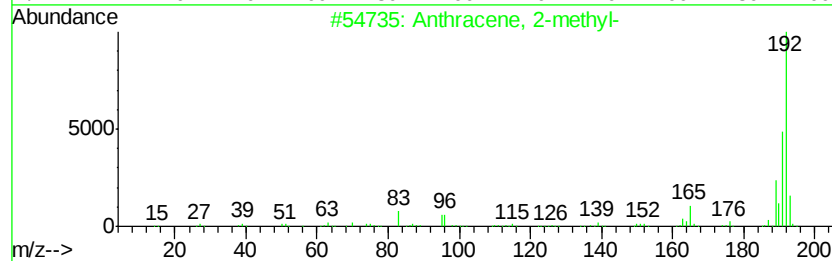
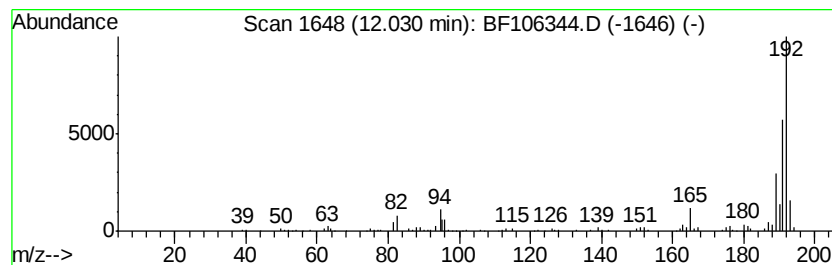
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Anthracene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.03	2.93 ng	147698	Phenanthrene-d10	11.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
2			Anthracene, 9-methyl-	192	C15H12	000779-02-2	96
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4			1H-Cyclopropa[1,1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	95
5			1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061218\
Data File : BF106344.D
Acq On : 12 Jun 2018 15:07
Operator : JU/SJ
Sample : J3245-03 2X
Misc :
ALS Vial : 9 Sample Multiplier: 1

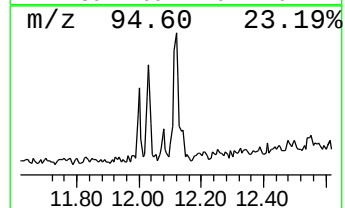
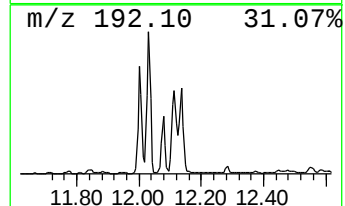
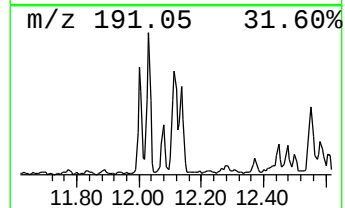
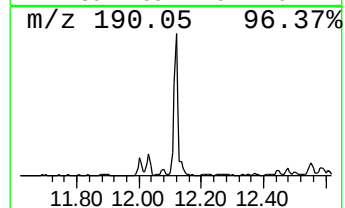
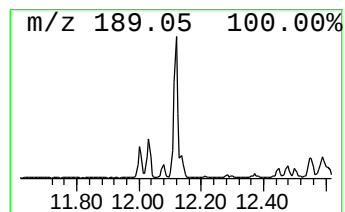
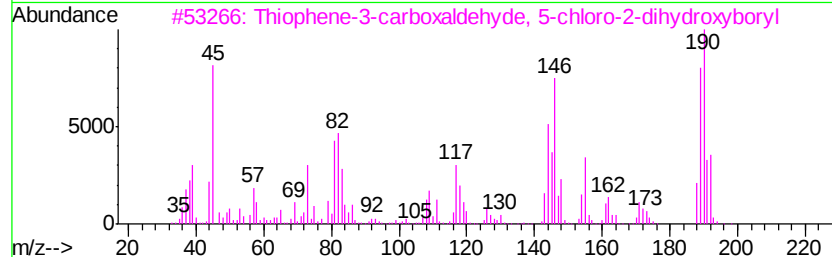
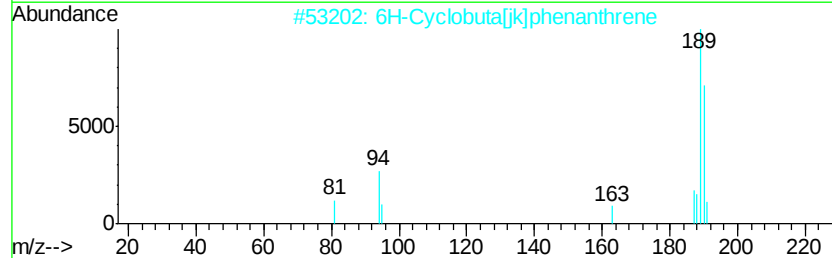
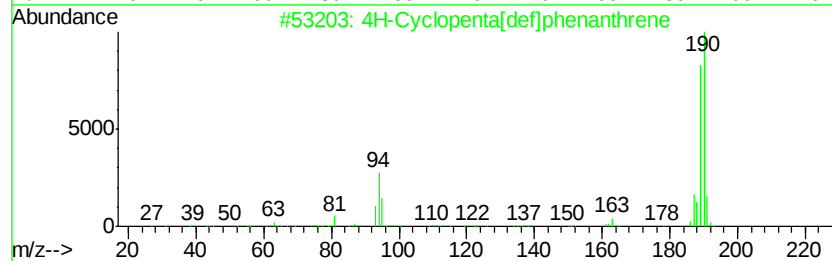
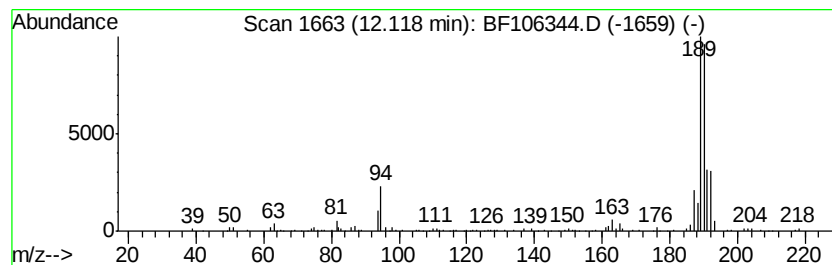
Instrument :
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ClientSampleId :
CL-01-061118-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.12	5.52 ng	277658	Phenanthrene-d10	11.50	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	94
2	6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	64
3	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	52
4	2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	47
5	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061218\
Data File : BF106344.D
Acq On : 12 Jun 2018 15:07
Operator : JU/SJ
Sample : J3245-03 2X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CL-01-061118-B

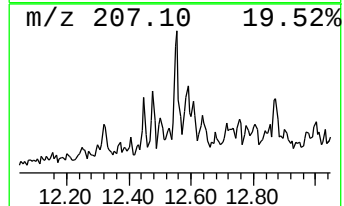
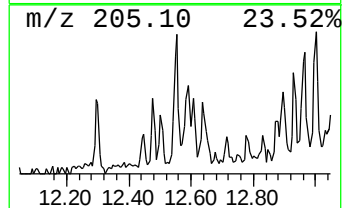
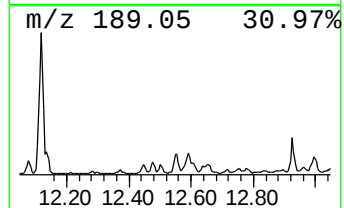
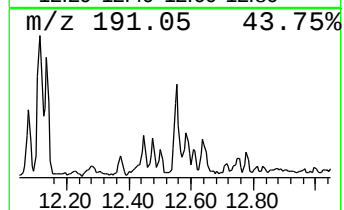
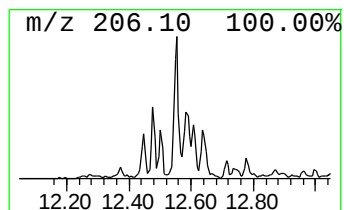
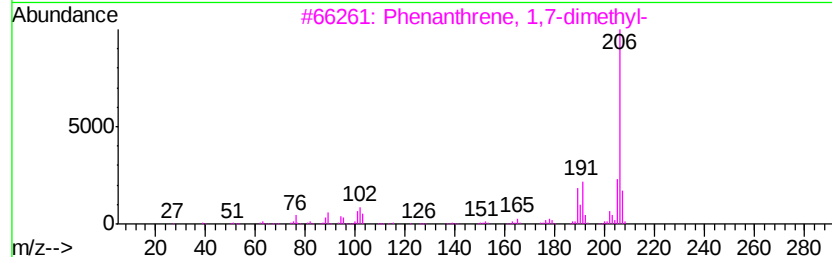
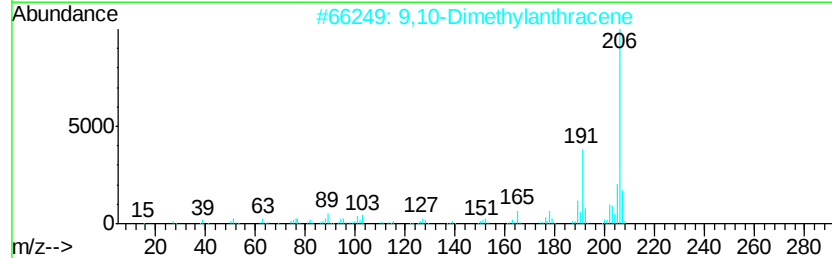
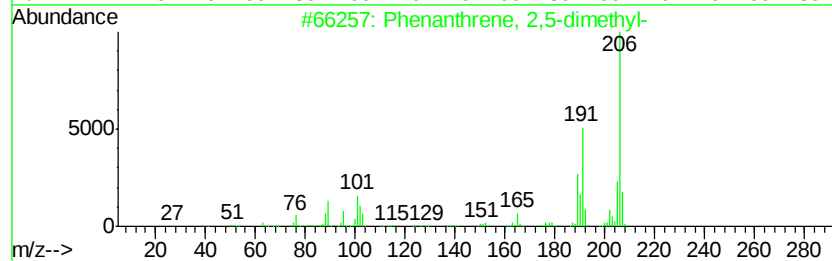
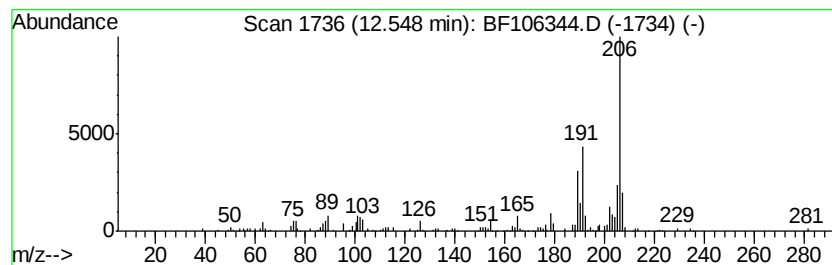
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 2,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.55	2.34 ng	117895	Phenanthrene-d10	11.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	97
2		9,10-Dimethylantracene	206	C16H14	000781-43-1	96
3		Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	93
4		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	90
5		di-p-Tolylacetylene	206	C16H14	002789-88-0	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061218\
Data File : BF106344.D
Acq On : 12 Jun 2018 15:07
Operator : JU/SJ
Sample : J3245-03 2X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CL-01-061118-B

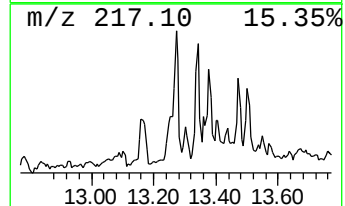
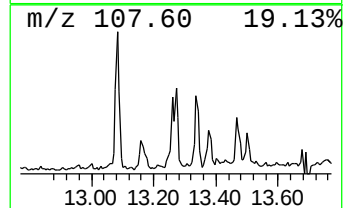
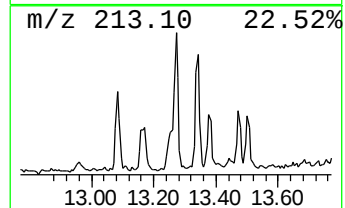
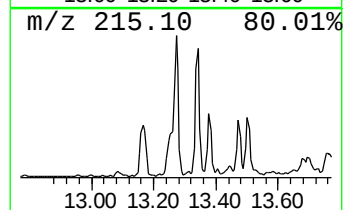
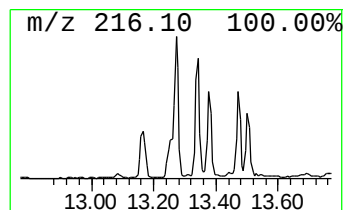
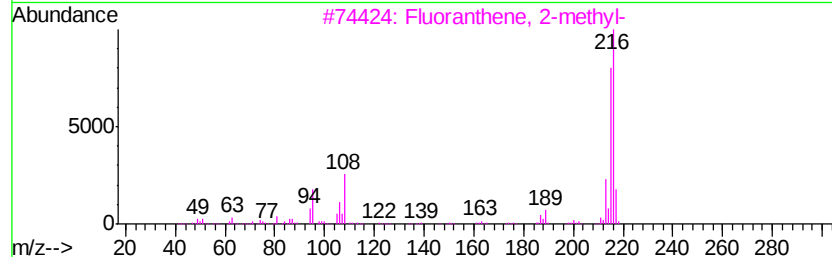
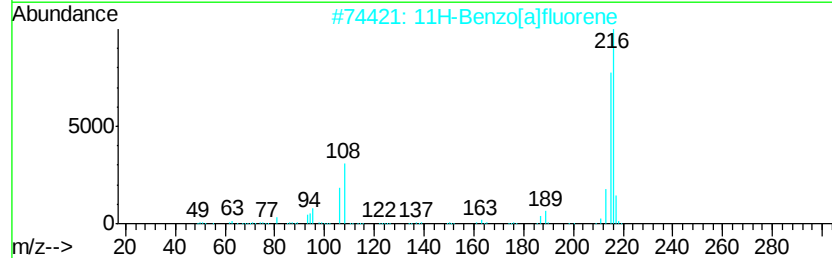
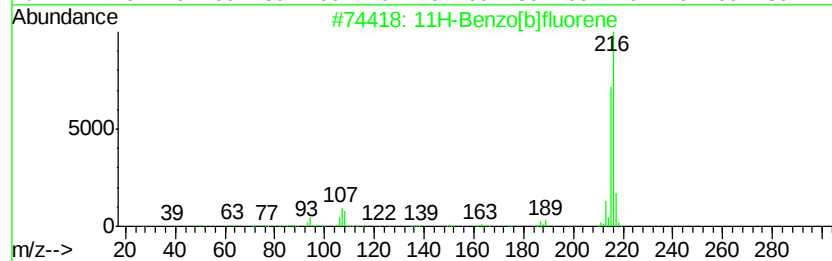
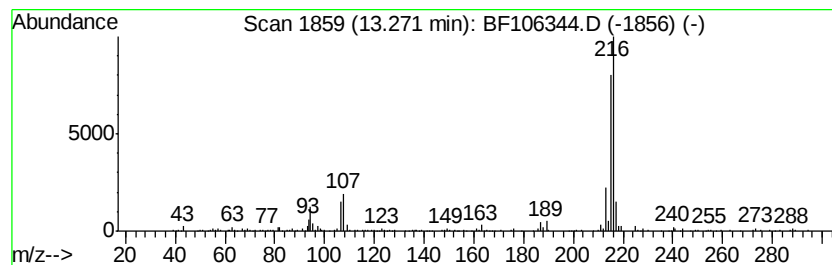
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.27	2.36 ng	225022	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
2		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81
3		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
4		Pvrene, 1-methyl-	216	C17H12	002381-21-7	64
5		1,3,5-Triazine-2,4,6-triamine, N...	216	C10H12N6	046731-79-7	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061218\
Data File : BF106344.D
Acq On : 12 Jun 2018 15:07
Operator : JU/SJ
Sample : J3245-03 2X
Misc :
ALS Vial : 9 Sample Multiplier: 1

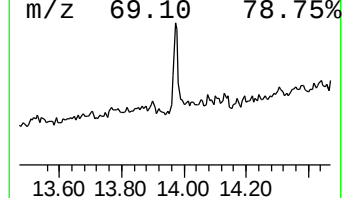
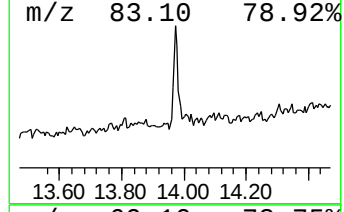
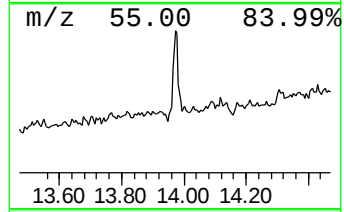
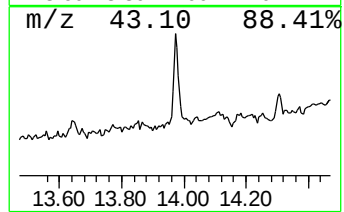
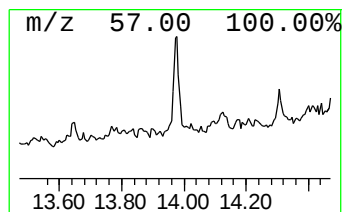
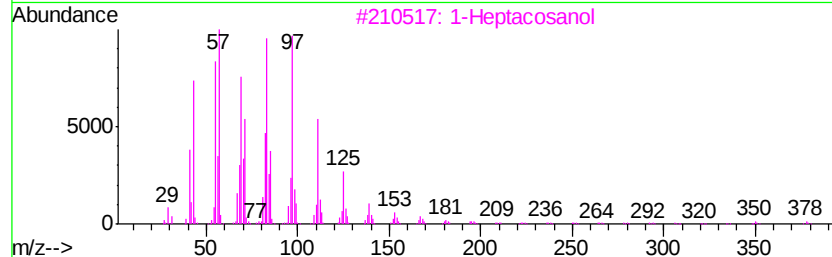
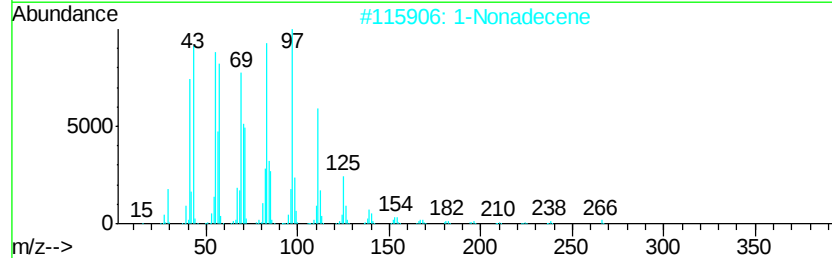
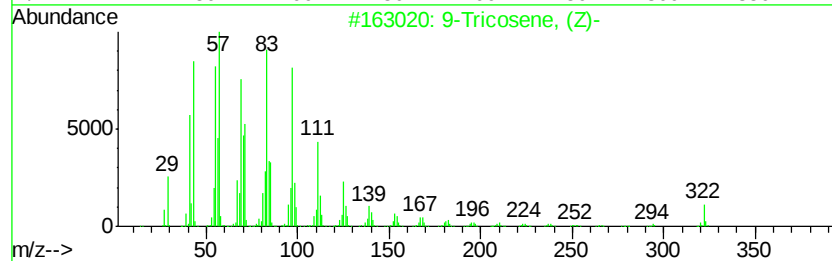
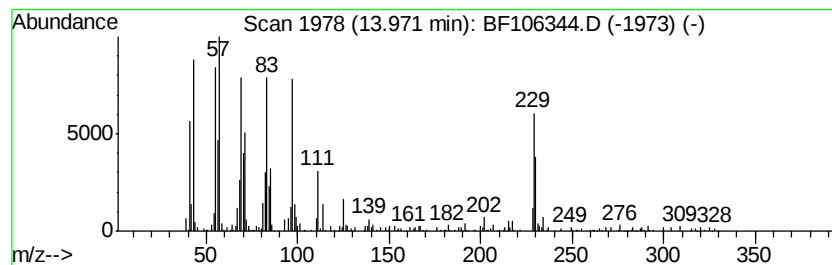
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 9-Tricosene, (Z)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.97	4.60 ng	437862	Chrysene-d12		14.16
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Tricosene, (Z)-	322	C23H46	027519-02-4	86
2	1-Nonadecene	266	C19H38	018435-45-5	76
3	1-Heptacosanol	396	C27H56O	002004-39-9	74
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	74
5	1-Hexadecanol	242	C16H34O	036653-82-4	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF061218\
Data File : BF106344.D
Acq On : 12 Jun 2018 15:07
Operator : JU/SJ
Sample : J3245-03 2X
Misc :
ALS Vial : 9 Sample Multiplier: 1

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	3.3	ng	150500	1	6.97	922612	20.0
unknown6.74	6.74	33.0	ng	1523610	1	6.97	922612	20.0
5H-Indeno[1,2-b]p...	11.73	2.2	ng	112525	4	11.50	1006920	20.0
Phenanthrene, 2-m...	12.00	3.3	ng	167958	4	11.50	1006920	20.0
Anthracene, 2-met...	12.03	2.9	ng	147698	4	11.50	1006920	20.0
4H-Cyclopenta[def...	12.12	5.5	ng	277658	4	11.50	1006920	20.0
Phenanthrene, 2,5...	12.55	2.3	ng	117895	4	11.50	1006920	20.0
11H-Benzo[b]fluorene	13.27	2.4	ng	225022	5	14.16	1905730	20.0
9-Tricosene, (Z)-	13.97	4.6	ng	437862	5	14.16	1905730	20.0