

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF080118\
 Data File : BF107891.D
 Acq On : 1 Aug 2018 16:13
 Operator : JU/SJ
 Sample : J4164-09 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-2(3-5)

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-BF073018.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	5.178	480	483	491	rBV	68396	89310	5.58%	0.415%
2	5.607	552	556	570	rBV	87065	303177	18.95%	1.409%
3	6.613	722	727	731	rBV	114138	240496	15.03%	1.118%
4	6.648	731	733	737	rVB	63304	75975	4.75%	0.353%
5	6.737	745	748	770	rVB	316201	560359	35.03%	2.605%
6	6.960	783	786	803	rBV	428403	777404	48.60%	3.614%
7	7.113	809	812	829	rBV	336818	486124	30.39%	2.260%
8	7.542	878	885	900	rBV2	68699	284597	17.79%	1.323%
9	8.242	1001	1004	1022	rBV	748789	1177798	73.63%	5.475%
10	9.054	1138	1142	1150	rBV	25589	38383	2.40%	0.178%
11	9.313	1183	1186	1200	rBV	422192	670299	41.90%	3.116%
12	9.589	1228	1233	1237	rBV3	15641	22751	1.42%	0.106%
13	9.660	1241	1245	1247	rBV	27589	33636	2.10%	0.156%
14	9.707	1247	1253	1262	rVB2	26035	53988	3.38%	0.251%
15	9.866	1276	1280	1285	rBV	40624	52378	3.27%	0.243%
16	10.001	1299	1303	1307	rBV	1039773	1067502	66.73%	4.962%
17	10.031	1307	1308	1318	rVB	141355	177683	11.11%	0.826%
18	10.201	1334	1337	1338	rBV	19101	19586	1.22%	0.091%
19	10.242	1342	1344	1348	rVB3	21903	21417	1.34%	0.100%
20	10.313	1353	1356	1359	rBV2	18174	22799	1.43%	0.106%
21	10.407	1368	1372	1378	rVB3	28119	33936	2.12%	0.158%
22	10.554	1392	1397	1403	rBV2	62124	95271	5.96%	0.443%
23	10.630	1406	1410	1414	rVB4	15839	28747	1.80%	0.134%
24	10.707	1421	1423	1428	rVB	21355	22187	1.39%	0.103%
25	10.801	1434	1439	1451	rBV	231153	438182	27.39%	2.037%
26	10.889	1451	1454	1464	rVB6	32783	68800	4.30%	0.320%
27	11.101	1486	1490	1493	rBV2	27401	36894	2.31%	0.172%
28	11.225	1506	1511	1513	rBV4	22608	27287	1.71%	0.127%
29	11.248	1513	1515	1521	rVB3	18238	21207	1.33%	0.099%
30	11.330	1527	1529	1532	rBV2	23097	23655	1.48%	0.110%
31	11.401	1538	1541	1548	rVB	34106	52769	3.30%	0.245%
32	11.501	1554	1558	1560	rBV	778378	952968	59.57%	4.430%
33	11.525	1560	1562	1568	rVV	786605	961354	60.10%	4.469%
34	11.577	1568	1571	1593	rVV	164360	447170	27.95%	2.079%

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Integrator: RTE

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.760	1596	1602	1604	rVV5	40083	74245	4.64%	0.345%
36	11.783	1604	1606	1611	rVV2	31652	43556	2.72%	0.202%
37	11.830	1611	1614	1621	rVV3	35056	58326	3.65%	0.271%
38	11.919	1625	1629	1633	rBV5	12484	19890	1.24%	0.092%
39	12.001	1640	1643	1646	rBV	99243	123560	7.72%	0.574%
40	12.030	1646	1648	1653	rVB2	99688	111824	6.99%	0.520%
41	12.077	1653	1656	1659	rBV	31151	31183	1.95%	0.145%
42	12.113	1659	1662	1665	rBV2	163923	199002	12.44%	0.925%
43	12.295	1690	1693	1697	rBV	61165	86893	5.43%	0.404%
44	12.472	1721	1723	1726	rBV	19774	20509	1.28%	0.095%
45	12.548	1733	1736	1739	rBV2	73630	82744	5.17%	0.385%
46	12.589	1740	1743	1745	rBV2	20003	16873	1.05%	0.078%
47	12.642	1748	1752	1761	rVB4	48535	99439	6.22%	0.462%
48	12.719	1761	1765	1768	rBV	1136633	1242309	77.66%	5.775%
49	12.819	1779	1782	1787	rVB	38120	46493	2.91%	0.216%
50	12.872	1788	1791	1797	rVB2	40907	66297	4.14%	0.308%
51	12.948	1797	1804	1810	rBV	1032375	1290644	80.68%	6.000%
52	12.995	1810	1812	1819	rVB3	65832	79280	4.96%	0.369%
53	13.077	1823	1826	1837	rBV	464353	581047	36.32%	2.701%
54	13.166	1837	1841	1847	rVB3	70014	119698	7.48%	0.556%
55	13.219	1847	1850	1852	rBV3	14590	17222	1.08%	0.080%
56	13.254	1852	1856	1857	rBV3	47293	60877	3.81%	0.283%
57	13.277	1857	1860	1865	rVB2	118671	137071	8.57%	0.637%
58	13.342	1868	1871	1874	rBV	91790	95792	5.99%	0.445%
59	13.383	1874	1878	1880	rBV2	56346	68303	4.27%	0.318%
60	13.471	1890	1893	1896	rBV	47187	56160	3.51%	0.261%
61	13.507	1896	1899	1908	rVB4	56713	96942	6.06%	0.451%
62	13.624	1917	1919	1923	rBV4	15021	20521	1.28%	0.095%
63	13.760	1938	1942	1944	rBV4	27617	42827	2.68%	0.199%
64	13.795	1944	1948	1954	rVV	49289	81417	5.09%	0.378%
65	13.901	1963	1966	1968	rBV	79286	87684	5.48%	0.408%
66	13.924	1968	1970	1973	rVV	96394	116387	7.28%	0.541%
67	13.954	1973	1975	1981	rVV4	103612	191469	11.97%	0.890%
68	14.001	1981	1983	1989	rVB2	51156	83550	5.22%	0.388%
69	14.148	2003	2008	2011	rBV2	1074196	1599624	100.00%	7.436%
70	14.177	2011	2013	2024	rVV	550463	810026	50.64%	3.765%
71	14.260	2024	2027	2033	rVB2	95237	131605	8.23%	0.612%

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Integrator: RTE
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Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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72	14.366	2043	2045	2047	rBV3	43063	44457	2.78%	0.207%
73	14.501	2066	2068	2069	rBV	34400	26940	1.68%	0.125%
74	14.542	2072	2075	2078	rVV	92791	126266	7.89%	0.587%
75	14.577	2079	2081	2084	rVB2	53538	59788	3.74%	0.278%

76	15.230	2187	2192	2203	rBV	440601	1138231	71.16%	5.291%
77	15.348	2208	2212	2216	rBV	81284	125211	7.83%	0.582%
78	15.554	2243	2247	2255	rBV	225423	361495	22.60%	1.680%
79	15.624	2255	2259	2265	rBV	368592	599532	37.48%	2.787%
80	15.689	2265	2270	2274	rBV	598593	880691	55.06%	4.094%

81	17.277	2534	2540	2550	rVB2	132150	335496	20.97%	1.560%
82	17.759	2614	2622	2632	rBV3	103774	336348	21.03%	1.564%

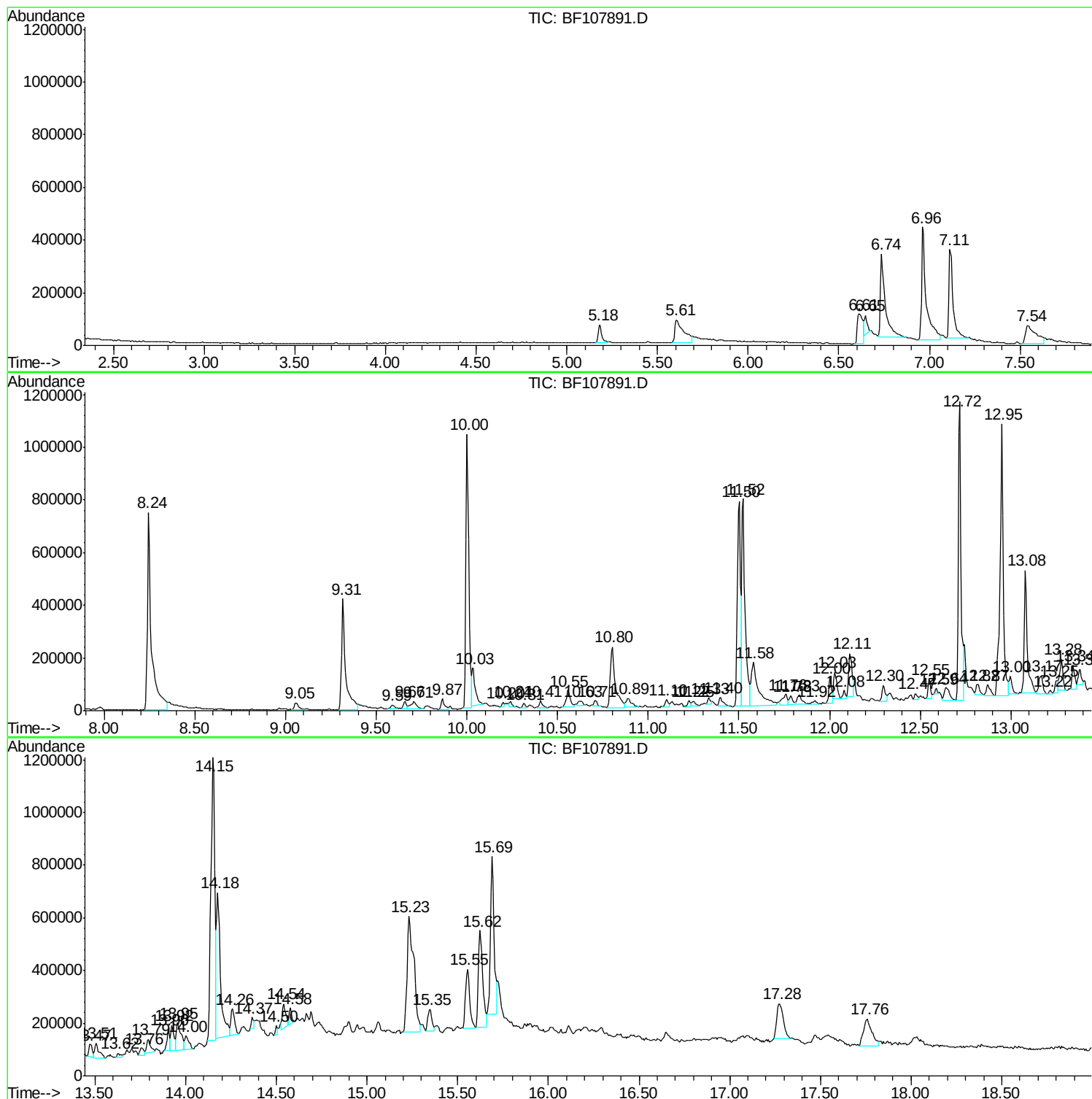
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
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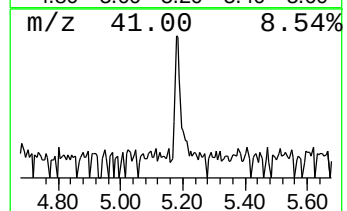
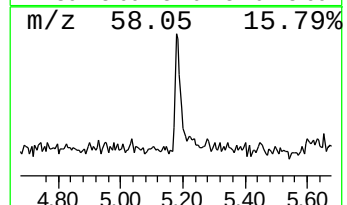
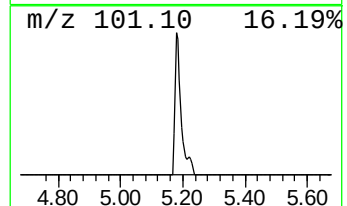
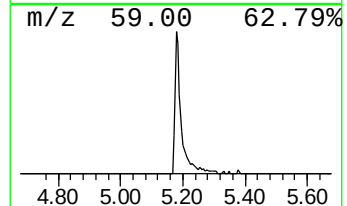
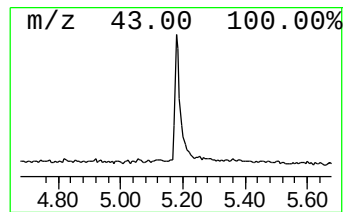
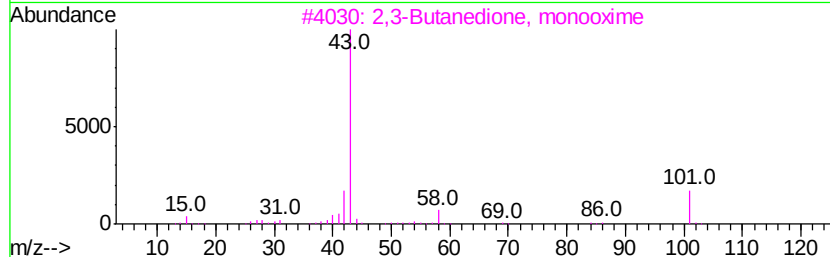
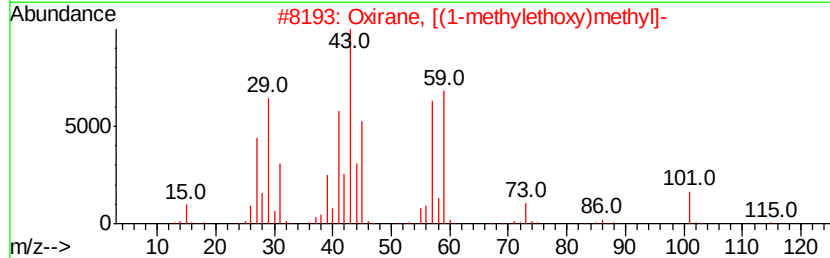
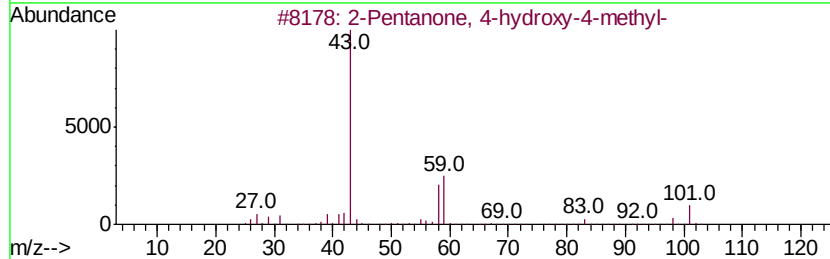
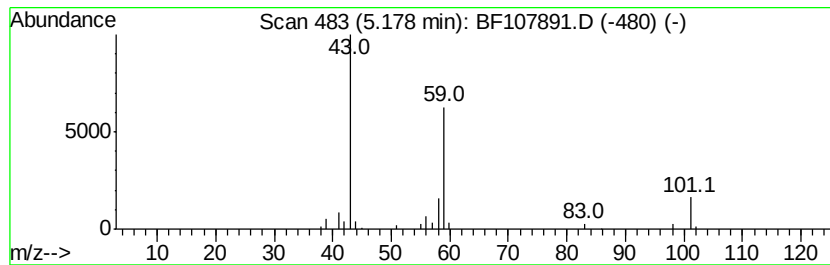
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	2.30 ng	89310	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	39
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
4		1,2-Butanediol	90	C4H10O2	000584-03-2	28
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28



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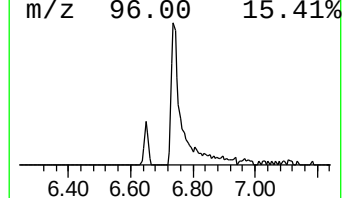
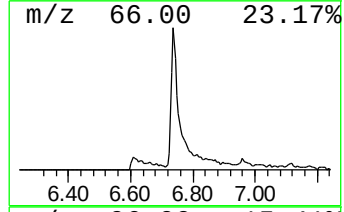
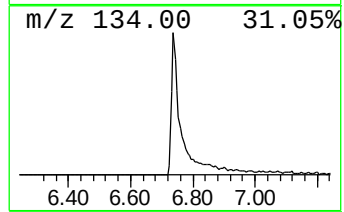
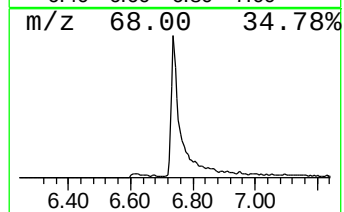
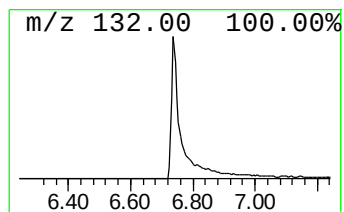
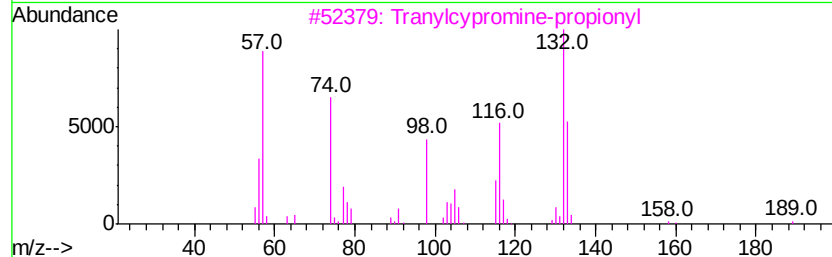
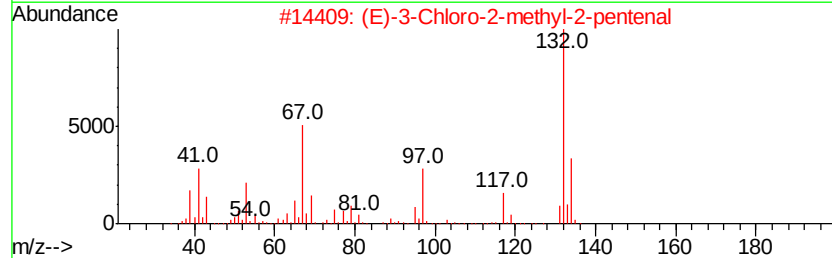
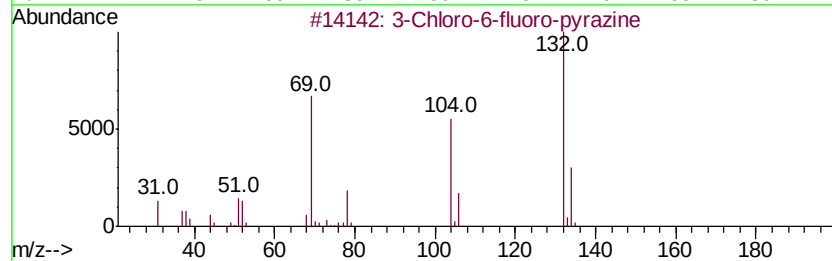
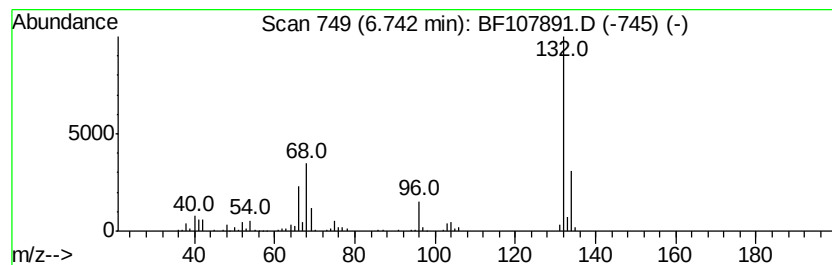
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	14.42 ng	560359	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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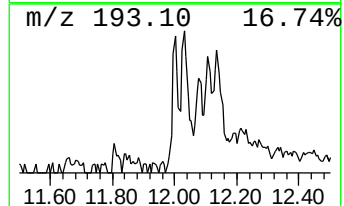
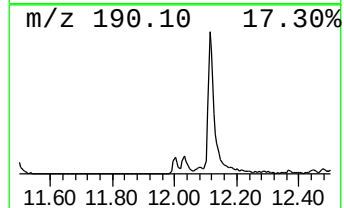
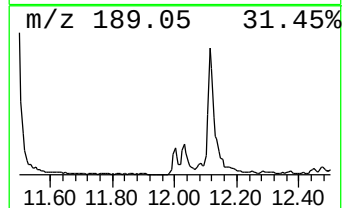
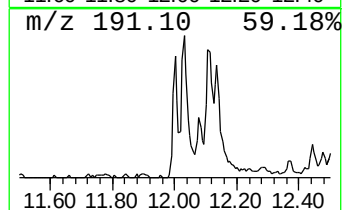
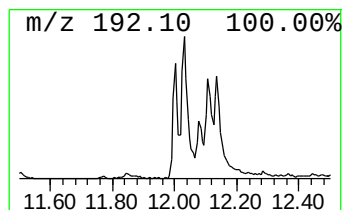
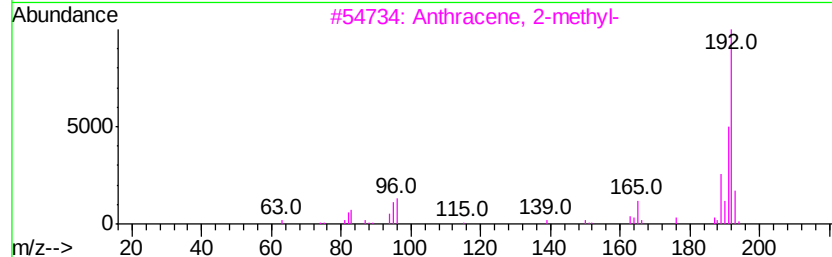
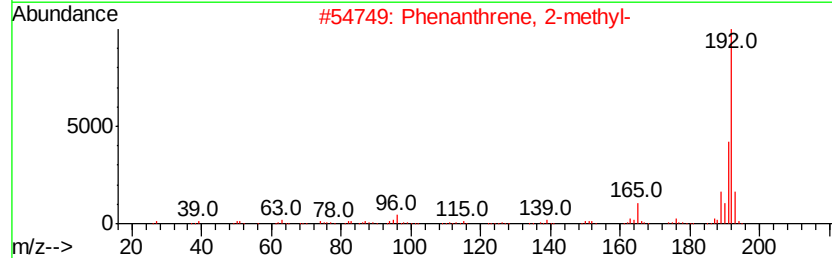
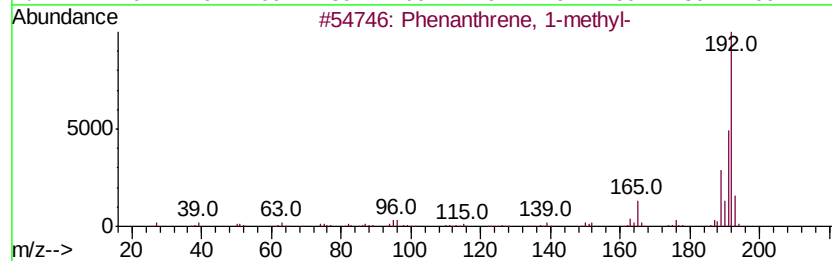
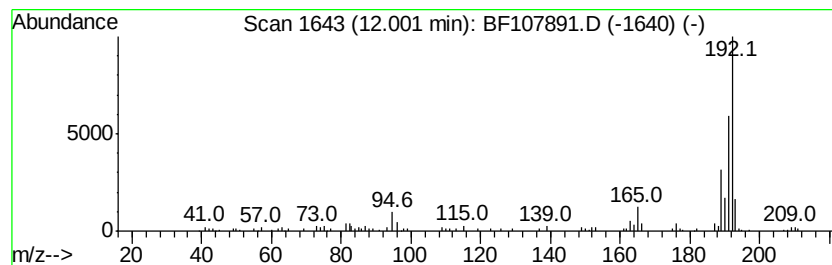
Instrument :
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ClientSampleId :
SB-2(3-5)

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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 4 Phenanthrene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.00	2.59 ng	123560	Phenanthrene-d10		11.50
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	93
5	1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	93



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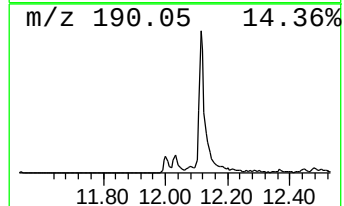
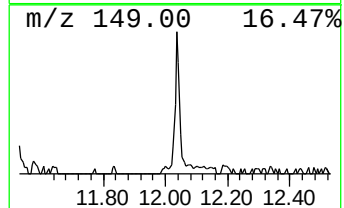
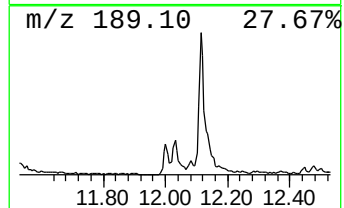
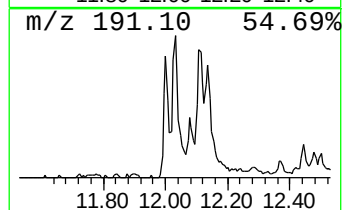
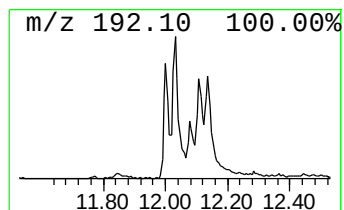
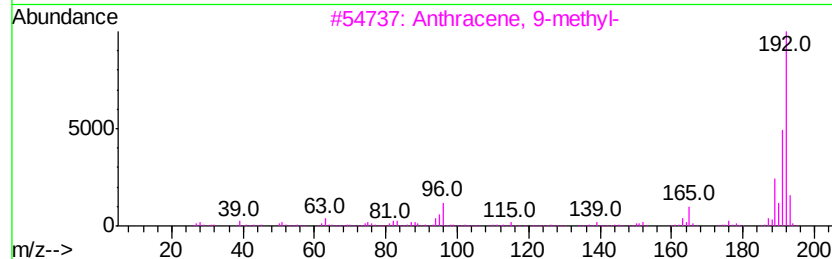
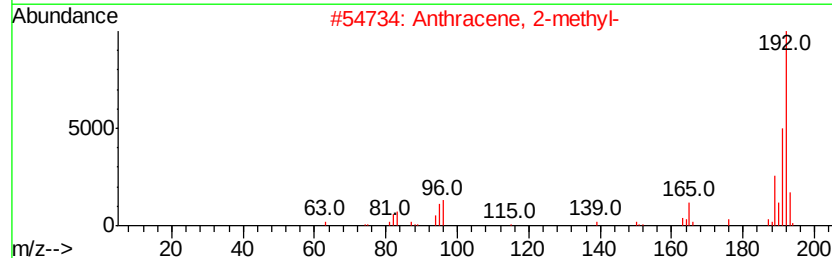
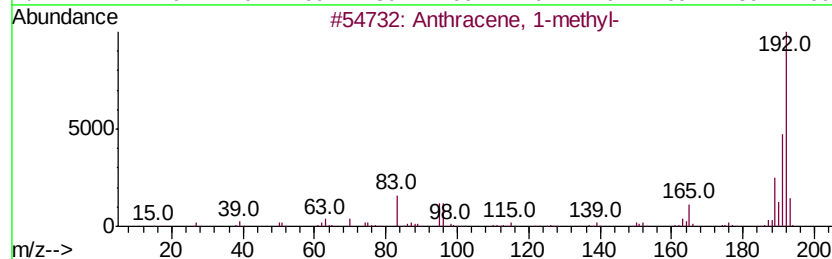
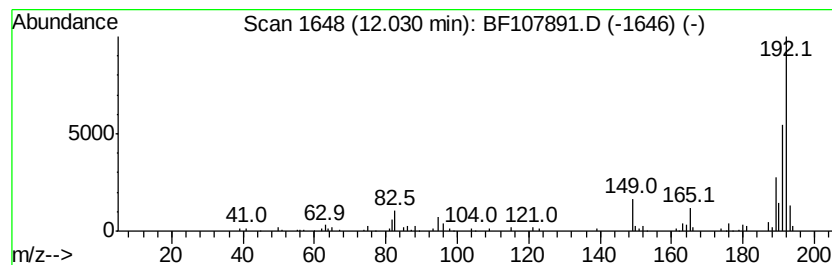
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Anthracene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.03	2.35 ng	111824	Phenanthrene-d10	11.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
3		Anthracene, 9-methyl-	192	C15H12	000779-02-2	89
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	83
5		1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080118\
Data File : BF107891.D
Acq On : 1 Aug 2018 16:13
Operator : JU/SJ
Sample : J4164-09 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

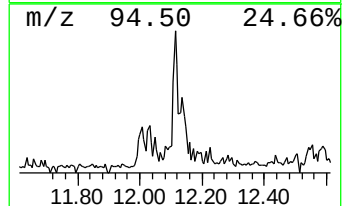
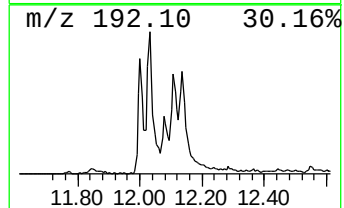
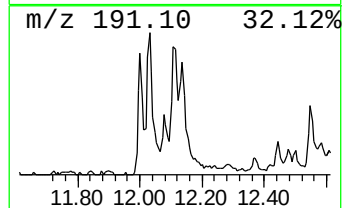
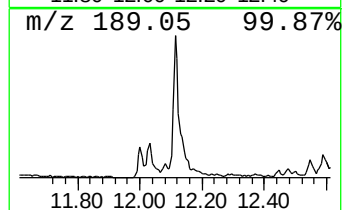
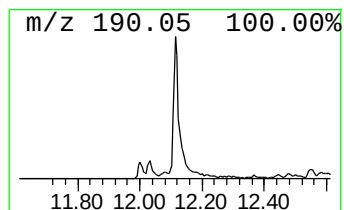
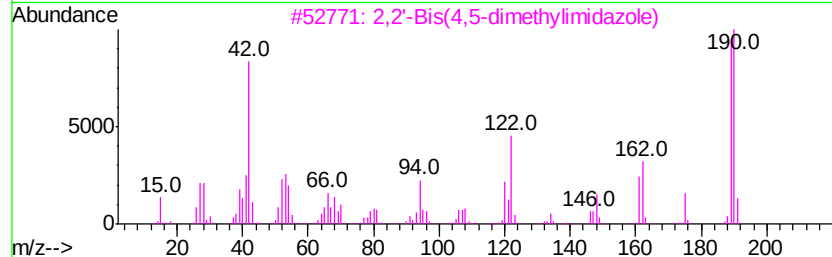
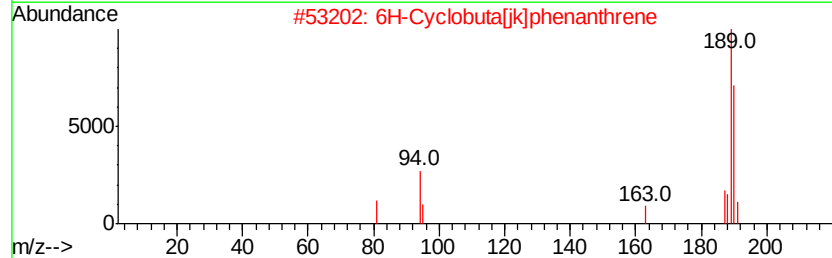
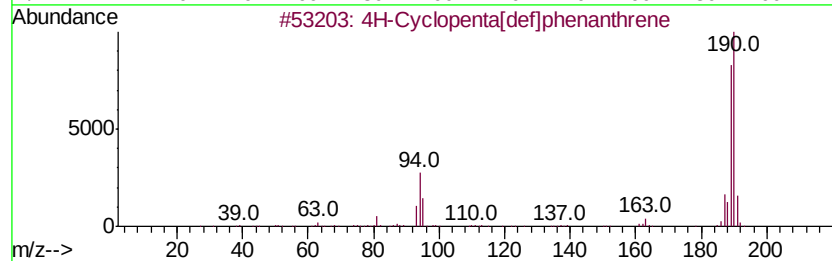
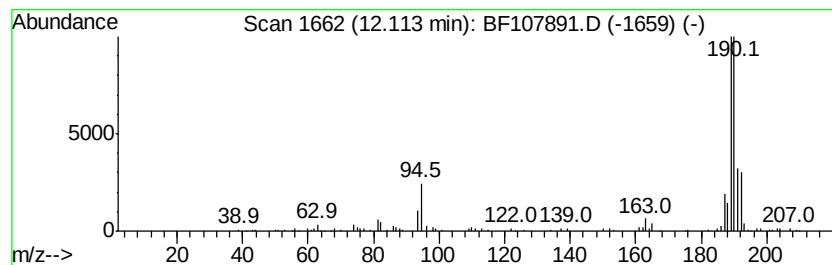
Instrument :
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ClientSampleId :
SB-2(3-5)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.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 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.11	4.18 ng	199002	Phenanthrene-d10	11.50	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	95
2	6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	72
3	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45
4	Methyl diselenide	190	C2H6Se2	007101-31-7	43
5	2,6-Dichlorobenzaldoxime	189	C7H5Cl2NO	025185-95-9	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080118\
Data File : BF107891.D
Acq On : 1 Aug 2018 16:13
Operator : JU/SJ
Sample : J4164-09 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-2(3-5)

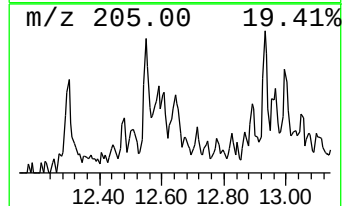
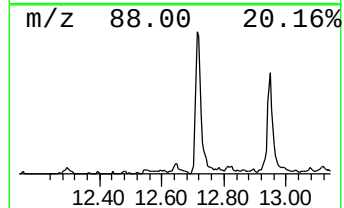
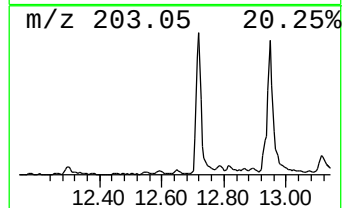
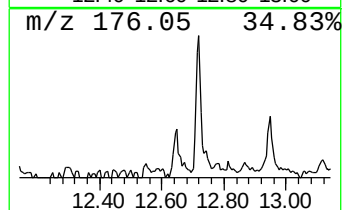
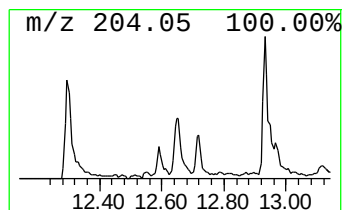
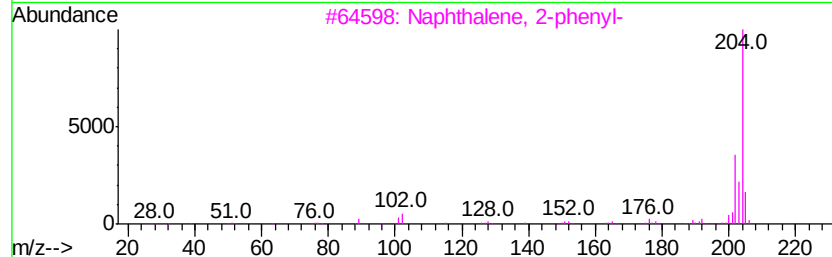
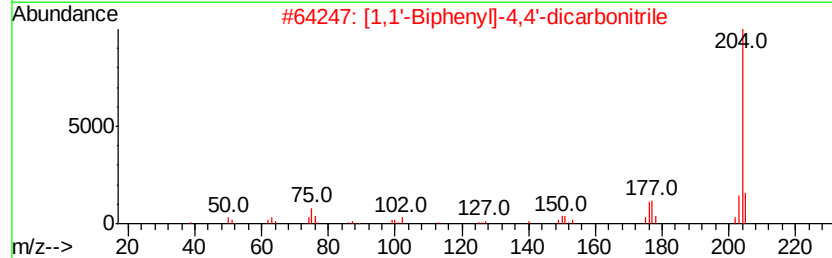
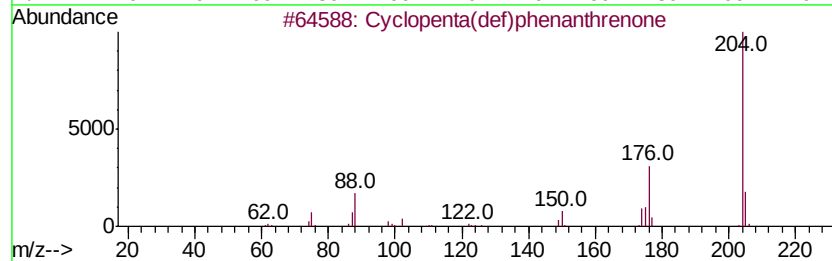
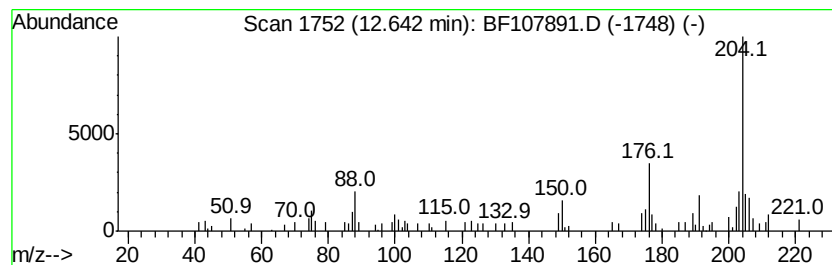
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Cyclopenta(def)phenanthrenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.64	2.09 ng	99439	Phenanthrene-d10	11.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	70
2		[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	70
3		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	53
4		9-Acridinecarbonitrile	204	C14H8N2	005326-19-2	50
5		[1,2,4]Triazolo[5,1-b]quinazolin...	204	C10H12N4O	1000337-56-5	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080118\
Data File : BF107891.D
Acq On : 1 Aug 2018 16:13
Operator : JU/SJ
Sample : J4164-09 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

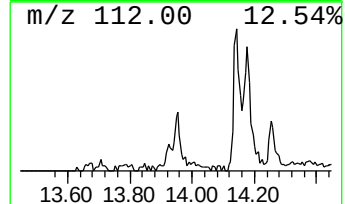
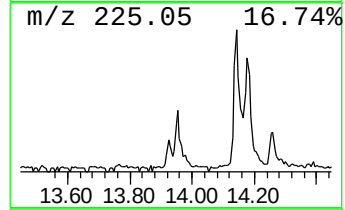
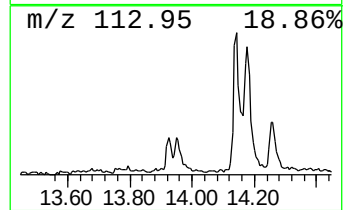
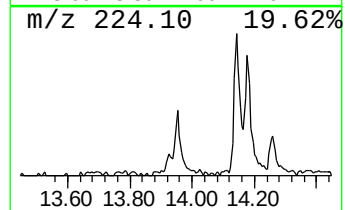
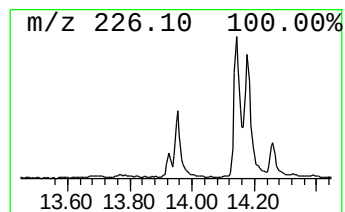
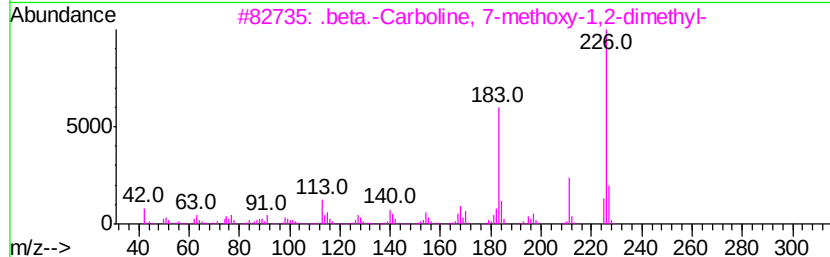
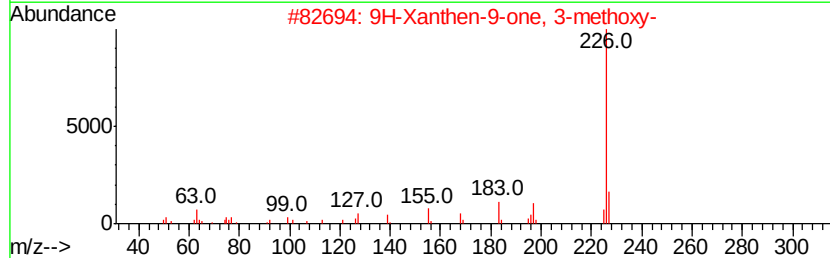
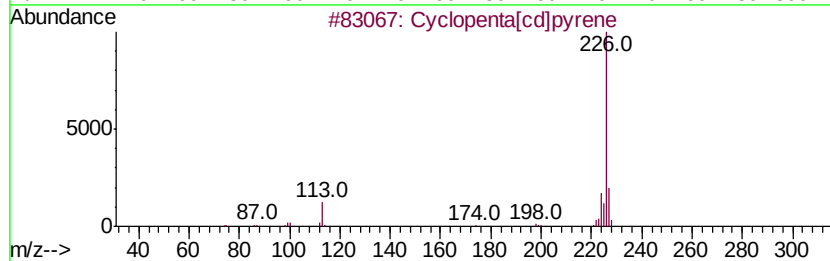
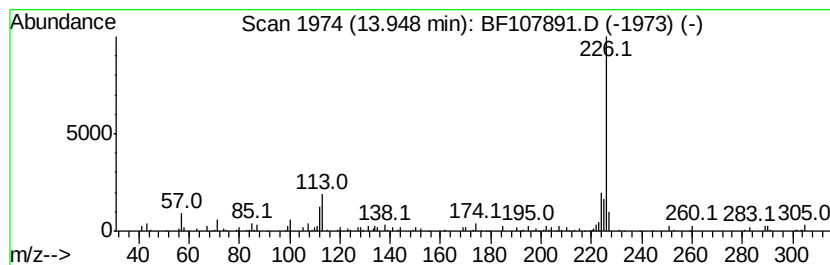
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Cyclopenta[cd]pyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.95	2.39 ng	191469	Chrysene-d12	14.15		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	52
2		9H-Xanthen-9-one, 3-methoxy-	226	C14H10O3	003722-52-9	47
3		.beta.-Carboline, 7-methoxy-1,2-...	226	C14H14N2O	006519-18-2	47
4		Benzo[ghi]fluoranthene	226	C18H10	000203-12-3	42
5		p-Anisaldehyde phenylhydrazone	226	C14H14N2O	000622-73-1	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080118\
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Acq On : 1 Aug 2018 16:13
Operator : JU/SJ
Sample : J4164-09 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-2(3-5)

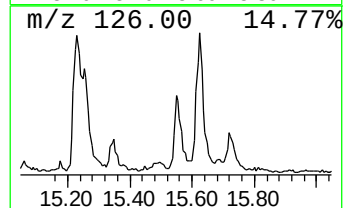
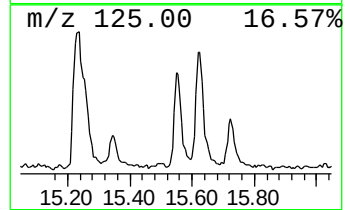
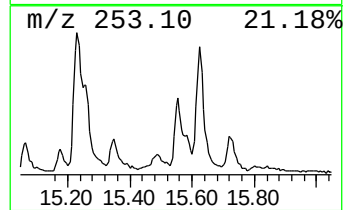
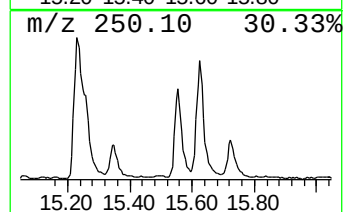
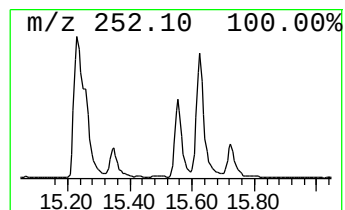
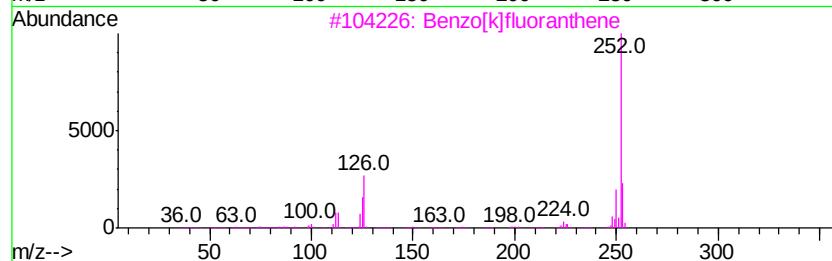
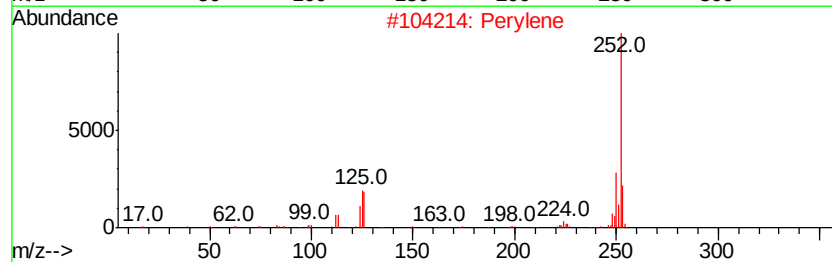
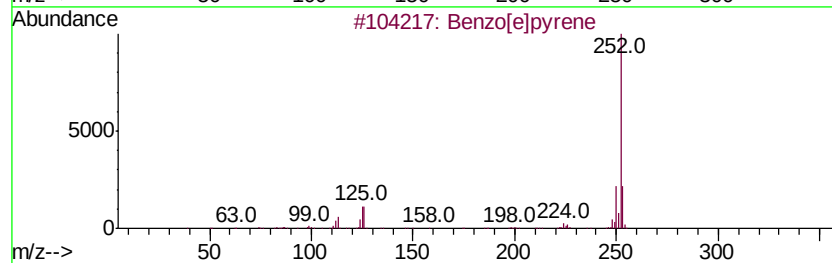
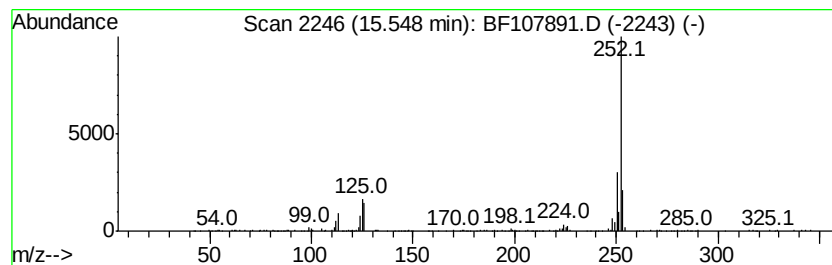
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	8.21 ng	361495	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	99
2		Perylene	252	C20H12	000198-55-0	94
3		Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
4		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	93
5		Benzo[j]fluoranthene	252	C20H12	000205-82-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF080118\
Data File : BF107891.D
Acq On : 1 Aug 2018 16:13
Operator : JU/SJ
Sample : J4164-09 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-2(3-5)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073018.M
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
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unknown6.74	6.74	14.4	ng	560359	1	6.96	777404	20.0
Phenanthrene, 1-m...	12.00	2.6	ng	123560	4	11.50	952968	20.0
Anthracene, 1-met...	12.03	2.4	ng	111824	4	11.50	952968	20.0
4H-Cyclopenta[def...	12.11	4.2	ng	199002	4	11.50	952968	20.0
Cyclopenta(def)ph...	12.64	2.1	ng	99439	4	11.50	952968	20.0
Cyclopenta[cd]pyrene	13.95	2.4	ng	191469	5	14.15	1599620	20.0
Benzo[e]pyrene	15.55	8.2	ng	361495	6	15.69	880691	20.0