

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF100720\
 Data File : BF121867.D
 Acq On : 7 Oct 2020 11:47
 Operator : JU/CG
 Sample : L4190-11 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PAV-B18-092920-0-10

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-BF091020.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.763	280	285	290	rVB	10270	16371	1.13%	0.080%
2	4.957	485	488	496	rVB	15008	20066	1.38%	0.098%
3	5.381	556	560	582	rBV	1037414	1146554	78.84%	5.600%
4	6.404	730	734	752	rBV	1004304	1203919	82.79%	5.881%
5	6.763	791	795	805	rVB	936642	792211	54.48%	3.870%
6	7.322	887	890	902	rBV	808609	818588	56.29%	3.998%
7	8.045	1009	1013	1016	rBV	1171030	1024806	70.47%	5.006%
8	8.757	1130	1134	1141	rBV	42421	42901	2.95%	0.210%
9	8.857	1147	1151	1159	rVB	55604	50612	3.48%	0.247%
10	9.122	1192	1196	1207	rBV	1553971	1454191	100.00%	7.103%
11	9.386	1237	1241	1247	rBV2	21716	31447	2.16%	0.154%
12	9.457	1249	1253	1255	rVV	53920	51615	3.55%	0.252%
13	9.486	1255	1258	1260	rVV	30873	31634	2.18%	0.155%
14	9.516	1260	1263	1270	rVV	119761	108548	7.46%	0.530%
15	9.575	1270	1273	1281	rVV	30498	45424	3.12%	0.222%
16	9.657	1285	1287	1291	rVB	20154	17500	1.20%	0.085%
17	9.798	1303	1311	1314	rBV	1418148	1218037	83.76%	5.949%
18	9.828	1314	1316	1322	rVV	209722	185806	12.78%	0.908%
19	9.904	1326	1329	1333	rVB3	11722	17290	1.19%	0.084%
20	9.951	1333	1337	1342	rBV2	19474	27154	1.87%	0.133%
21	10.004	1342	1346	1350	rVV	106264	118987	8.18%	0.581%
22	10.039	1350	1352	1358	rVB2	26276	28736	1.98%	0.140%
23	10.116	1362	1365	1368	rBV2	20058	20870	1.44%	0.102%
24	10.204	1376	1380	1381	rBV2	22669	22754	1.56%	0.111%
25	10.228	1381	1384	1387	rVB2	33520	30263	2.08%	0.148%
26	10.345	1400	1404	1410	rBV2	149325	147657	10.15%	0.721%
27	10.392	1410	1412	1413	rBV	20967	15475	1.06%	0.076%
28	10.433	1417	1419	1421	rVV	30943	28953	1.99%	0.141%
29	10.480	1424	1427	1428	rVV	31149	35004	2.41%	0.171%
30	10.504	1428	1431	1434	rVB	60379	64217	4.42%	0.314%
31	10.586	1441	1445	1451	rBV	821290	757627	52.10%	3.701%
32	10.710	1462	1466	1469	rBV3	28145	42979	2.96%	0.210%
33	10.892	1491	1497	1500	rBV4	43616	62441	4.29%	0.305%
34	10.922	1500	1502	1505	rVB2	19934	15802	1.09%	0.077%

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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-BF091020.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	10.975	1509	1511	1514	rVB	18888	16847	1.16%	0.082%
36	11.022	1514	1519	1521	rBV2	36189	52380	3.60%	0.256%
37	11.092	1528	1531	1534	rVB	50146	44555	3.06%	0.218%
38	11.145	1534	1540	1543	rVV3	43606	76316	5.25%	0.373%
39	11.180	1543	1546	1553	rVB	85082	85455	5.88%	0.417%
40	11.280	1559	1563	1566	rBV	1206020	1296643	89.17%	6.333%
41	11.310	1566	1568	1571	rVV	1076276	863388	59.37%	4.217%
42	11.357	1571	1576	1579	rVV	300734	290083	19.95%	1.417%
43	11.398	1579	1583	1585	rVB2	28528	30264	2.08%	0.148%
44	11.516	1597	1603	1610	rBV2	98746	150631	10.36%	0.736%
45	11.575	1610	1613	1618	rVV2	46337	50115	3.45%	0.245%
46	11.627	1618	1622	1625	rVB5	16326	24410	1.68%	0.119%
47	11.675	1625	1630	1632	rBV5	14996	22620	1.56%	0.110%
48	11.745	1638	1642	1645	rBV4	13793	19756	1.36%	0.096%
49	11.786	1645	1649	1651	rBV	107901	113949	7.84%	0.557%
50	11.816	1651	1654	1658	rBV2	246293	227988	15.68%	1.114%
51	11.863	1659	1662	1664	rBV	58089	52914	3.64%	0.258%
52	11.898	1664	1668	1670	rBV	145674	156045	10.73%	0.762%
53	11.980	1677	1682	1687	rBV3	35418	66669	4.58%	0.326%
54	12.080	1696	1699	1701	rBV	83935	67146	4.62%	0.328%
55	12.110	1701	1704	1707	rBV	43778	36891	2.54%	0.180%
56	12.227	1722	1724	1727	rVB2	30228	26218	1.80%	0.128%
57	12.263	1727	1730	1731	rBV2	31925	27846	1.91%	0.136%
58	12.280	1731	1733	1736	rVB2	25310	24324	1.67%	0.119%
59	12.333	1739	1742	1744	rBV	63139	67230	4.62%	0.328%
60	12.369	1744	1748	1751	rVV3	58119	80432	5.53%	0.393%
61	12.422	1754	1757	1761	rVB3	41401	45153	3.11%	0.221%
62	12.457	1761	1763	1766	rBV3	15356	15876	1.09%	0.078%
63	12.498	1766	1770	1773	rBV	851937	764742	52.59%	3.735%
64	12.563	1778	1781	1783	rBV	23031	18733	1.29%	0.092%
65	12.610	1784	1789	1792	rBV	279572	317147	21.81%	1.549%
66	12.722	1802	1808	1814	rVV2	661374	800973	55.08%	3.912%
67	12.774	1815	1817	1822	rVB2	48394	63582	4.37%	0.311%
68	12.869	1827	1833	1836	rBV	1377183	1337787	92.00%	6.534%
69	13.051	1857	1864	1868	rBV2	105965	202108	13.90%	0.987%
70	13.121	1873	1876	1878	rVB	47952	46891	3.22%	0.229%
71	13.157	1878	1882	1885	rBV2	79071	103127	7.09%	0.504%

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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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	13.280	1901	1903	1906	rBV2	37437	38664	2.66%	0.189%
73	13.345	1912	1914	1922	rVB	130208	140773	9.68%	0.688%
74	13.563	1949	1951	1957	rVB	38184	42096	2.89%	0.206%
75	13.674	1967	1970	1972	rBV	43691	48118	3.31%	0.235%
76	13.698	1972	1974	1976	rBV	49929	45503	3.13%	0.222%
77	13.921	2007	2012	2015	rBV2	985976	1116365	76.77%	5.453%
78	13.945	2015	2016	2025	rVB	233437	195328	13.43%	0.954%
79	14.951	2184	2187	2194	rVB	186132	286589	19.71%	1.400%
80	15.304	2244	2247	2253	rBV	194800	256981	17.67%	1.255%
81	15.368	2254	2258	2265	rVB	674543	918939	63.19%	4.489%

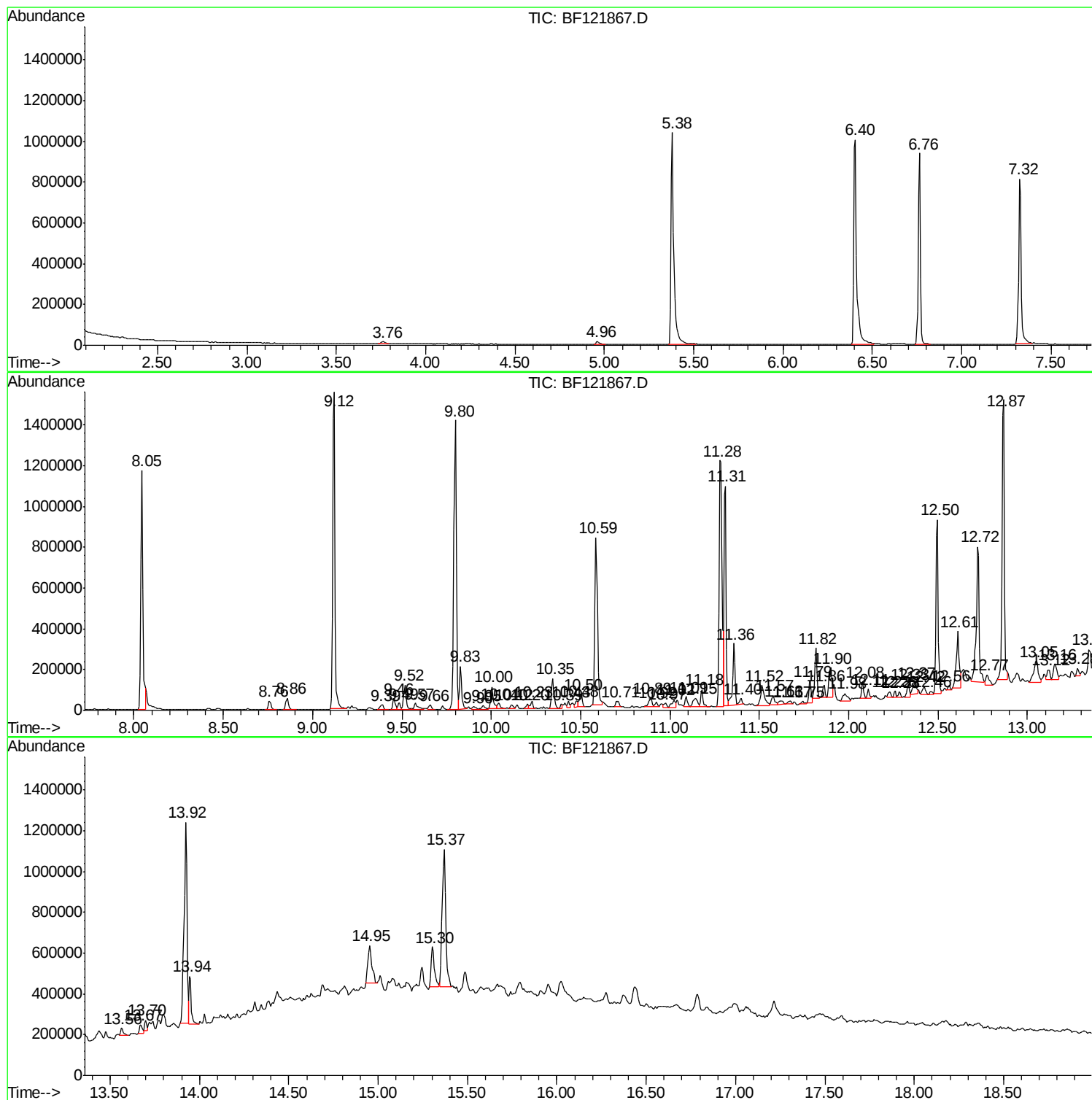
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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ALS Vial : 5 Sample Multiplier: 1

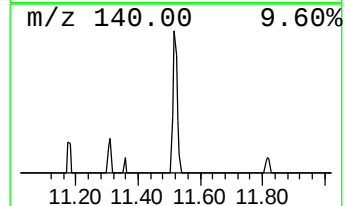
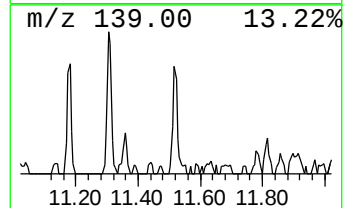
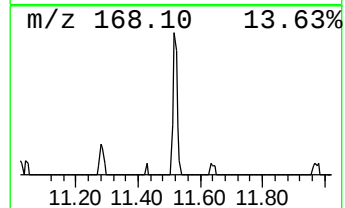
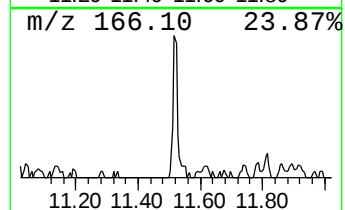
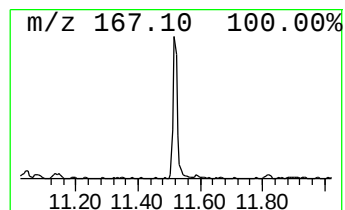
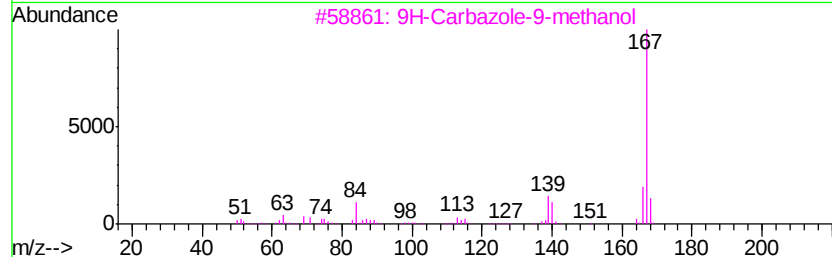
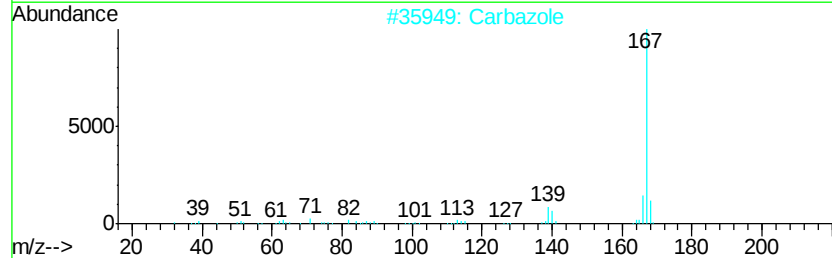
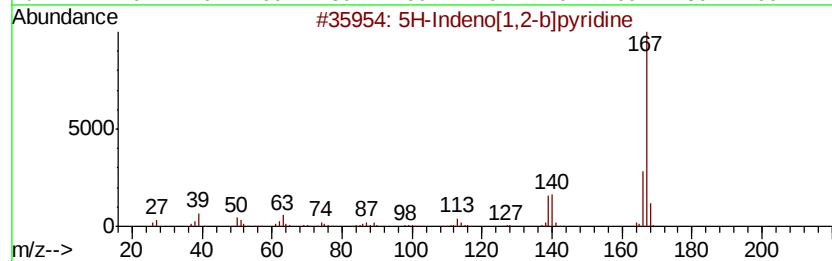
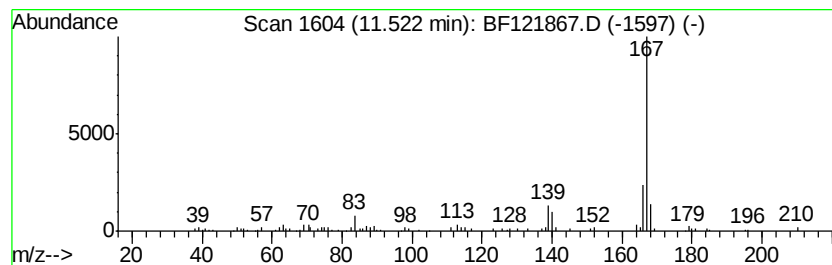
Instrument :
BNA_F
ClientSampleId :
PAV-B18-092920-0-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 5H-Indeno[1,2-b]pyridine Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.52	2.32 ng	150631	Phenanthrene-d10	11.29		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-Indeno[1,2-b]pvridine	167	C12H9N	000244-99-5	94	
2	Carbazole	167	C12H9N	000086-74-8	87	
3	9H-Carbazole-9-methanol	197	C13H11NO	002409-36-1	83	
4	9H-Carbazole, 9-nitroso-	196	C12H8N2O	002788-23-0	83	
5	D-Galactitol, 2-(acetylmethylami...	363	C16H29NO8	074410-42-7	80	



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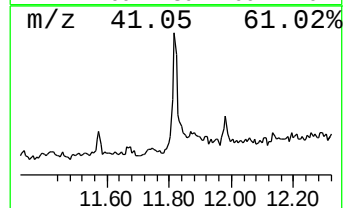
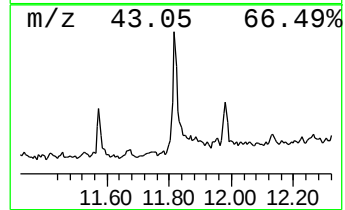
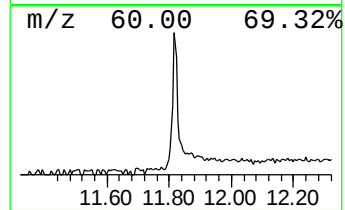
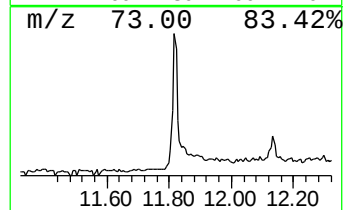
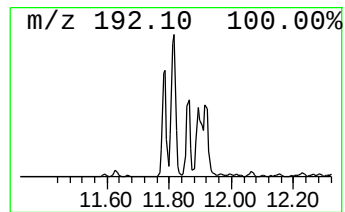
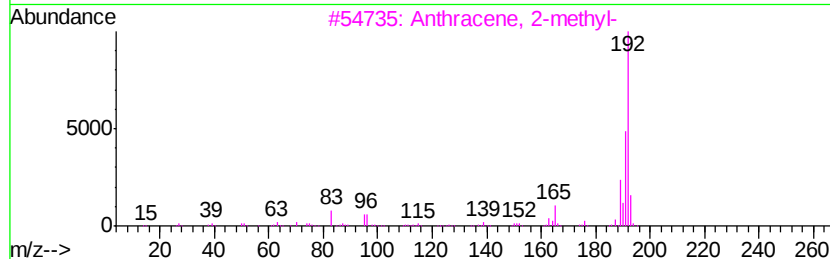
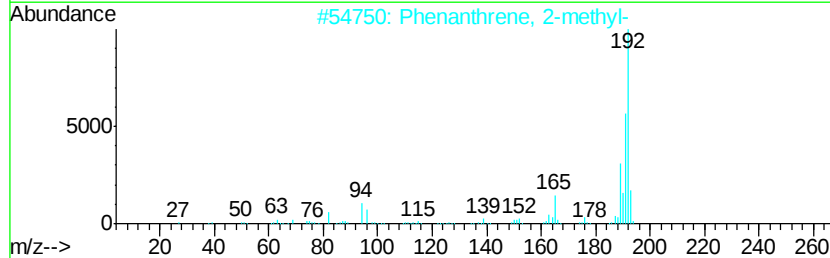
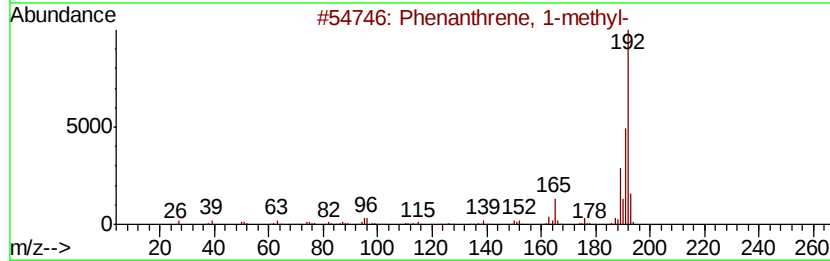
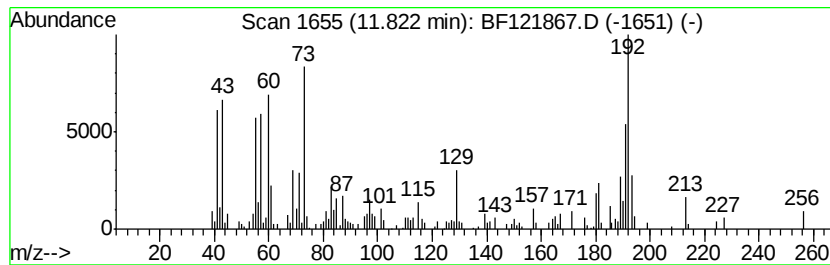
Instrument :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Phenanthrene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.82	3.52 ng	227988	Phenanthrene-d10		11.29
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	97
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	92
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	91



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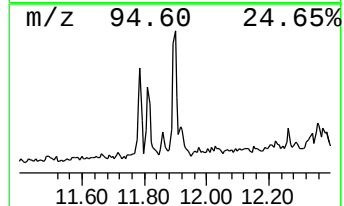
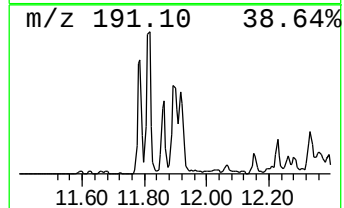
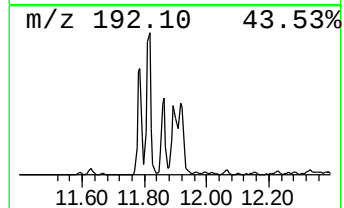
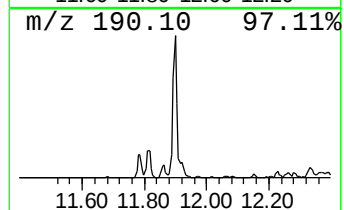
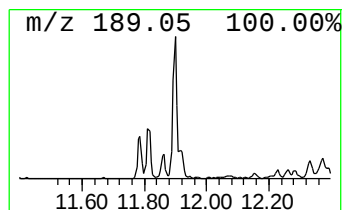
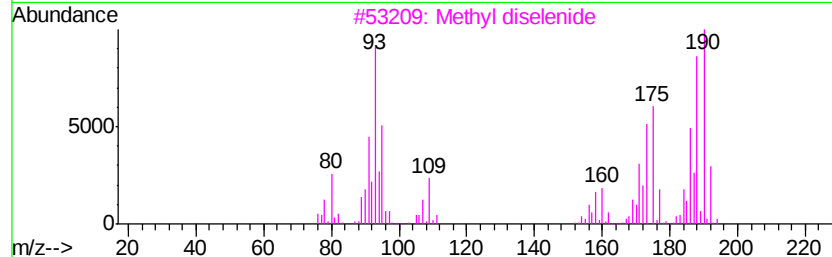
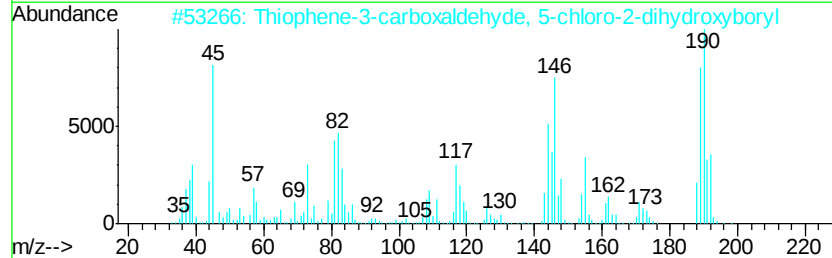
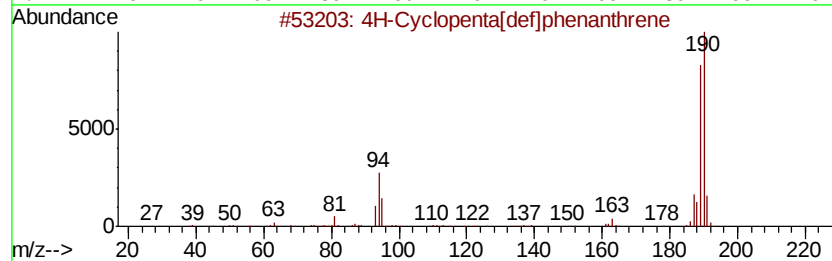
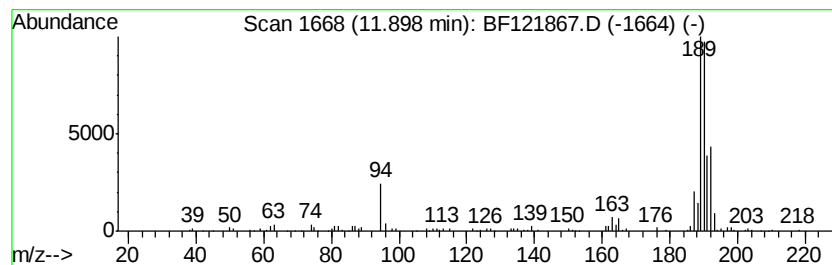
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.90	2.41 ng	156045	Phenanthrene-d10		11.29	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	50
3		Methyl diselenide	190	C2H6Se2	007101-31-7	45
4		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	43
5		6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	40



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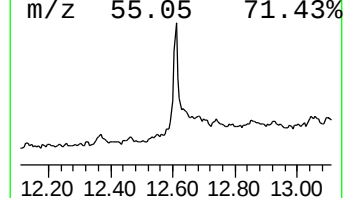
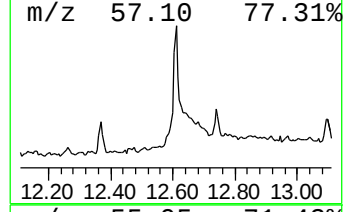
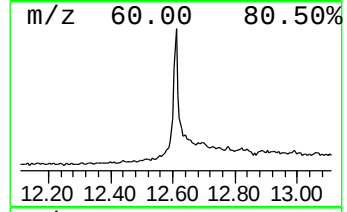
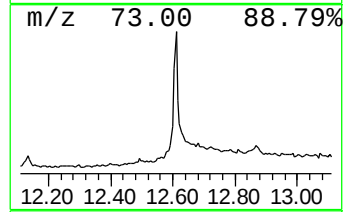
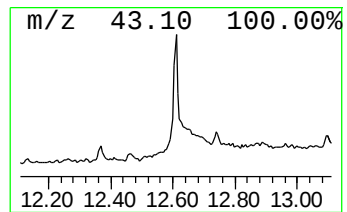
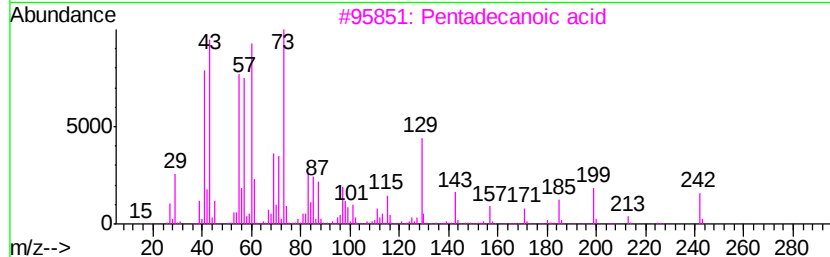
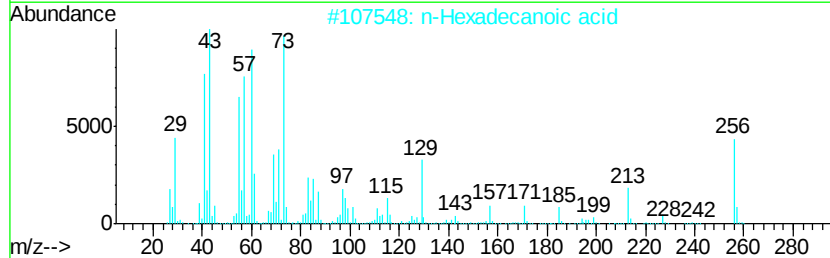
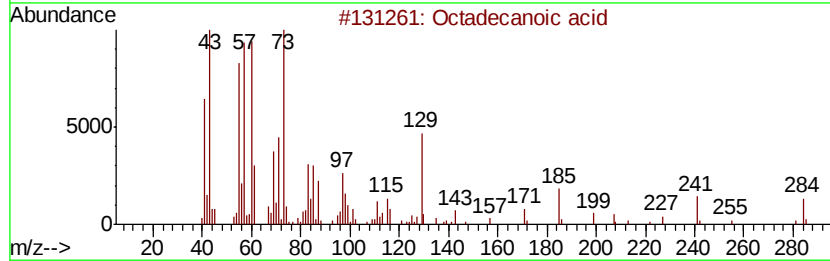
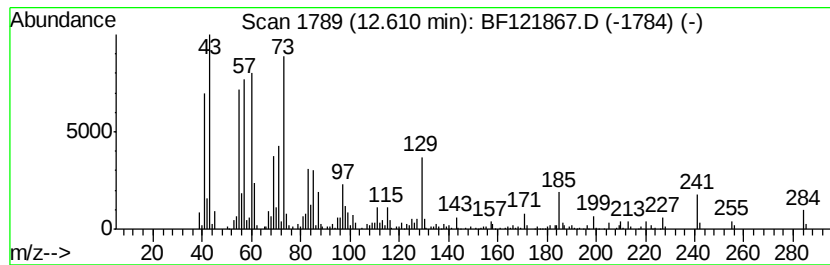
Instrument :
BNA_F
ClientSampleId :
PAV-B18-092920-0-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.61	5.68 ng	317147	Chrysene-d12		13.92
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4	Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	86
5	Tridecanoic acid	214	C13H26O2	000638-53-9	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100720\
Data File : BF121867.D
Acq On : 7 Oct 2020 11:47
Operator : JU/CG
Sample : L4190-11 2X
Misc :
ALS Vial : 5 Sample Multiplier: 1

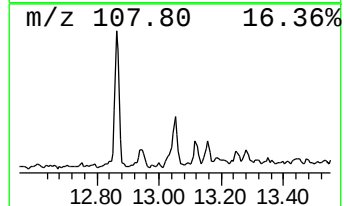
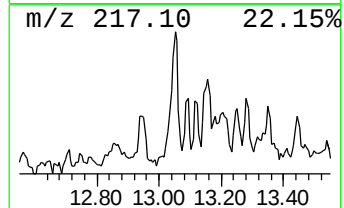
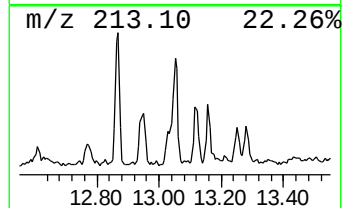
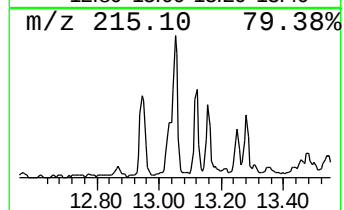
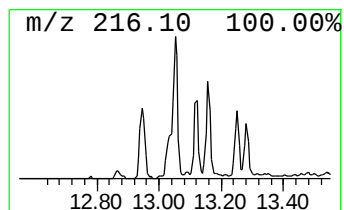
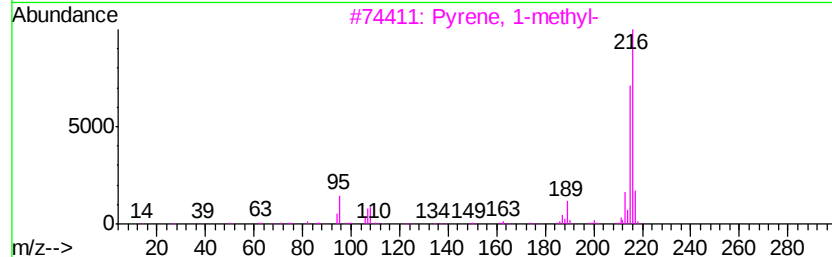
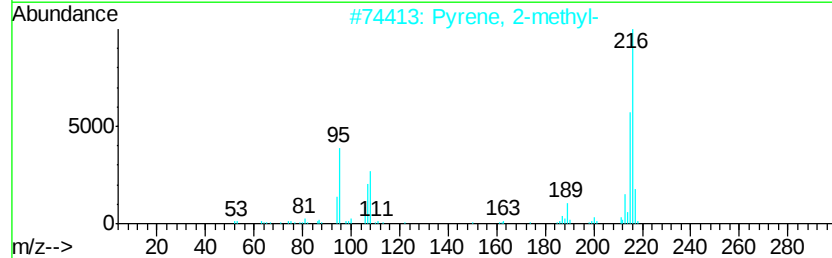
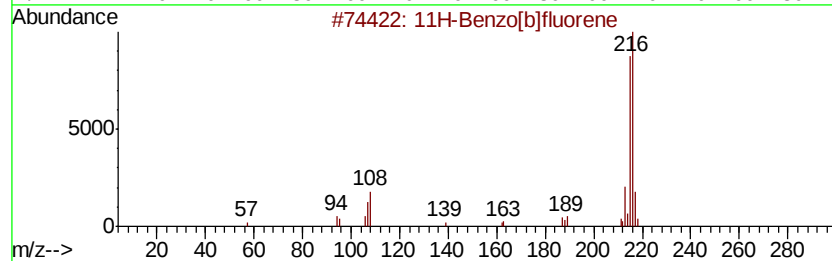
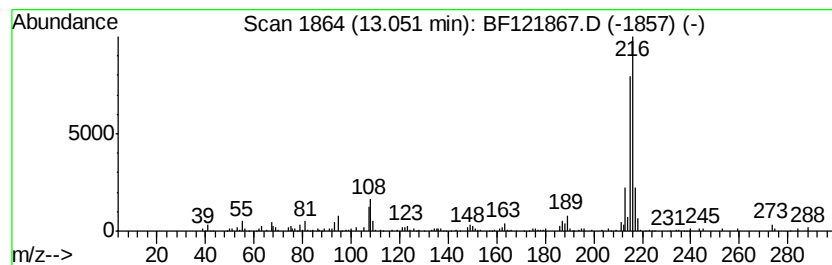
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ClientSampleId :
PAV-B18-092920-0-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 11H-Benzo[b]fluorene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.05	3.62 ng	202108	Chrysene-d12	13.92
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 96
2		Pyrene, 2-methyl-	216 C17H12	003442-78-2 94
3		Pyrene, 1-methyl-	216 C17H12	002381-21-7 94
4		Fluoranthene, 2-methyl-	216 C17H12	033543-31-6 93
5		11H-Benzo[a]fluorene	216 C17H12	000238-84-6 91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100720\
Data File : BF121867.D
Acq On : 7 Oct 2020 11:47
Operator : JU/CG
Sample : L4190-11 2X
Misc :
ALS Vial : 5 Sample Multiplier: 1

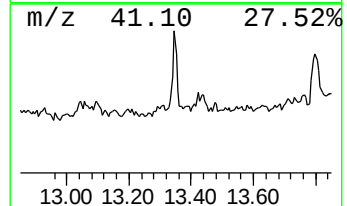
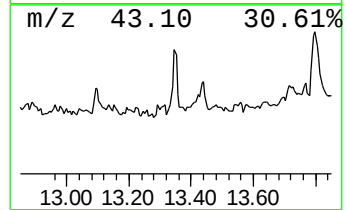
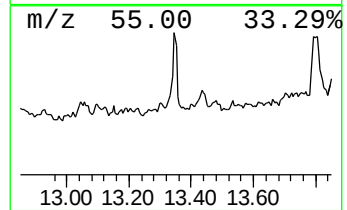
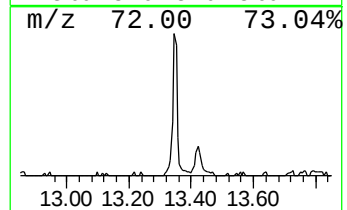
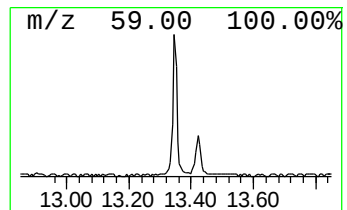
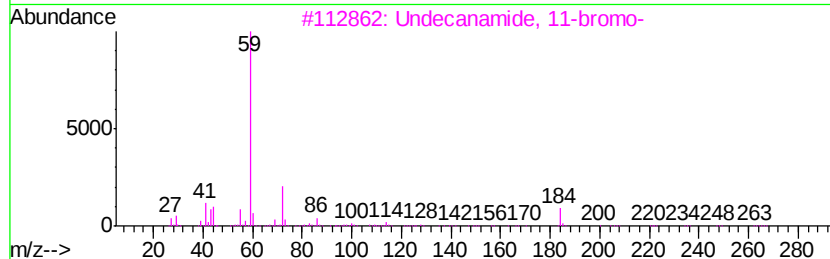
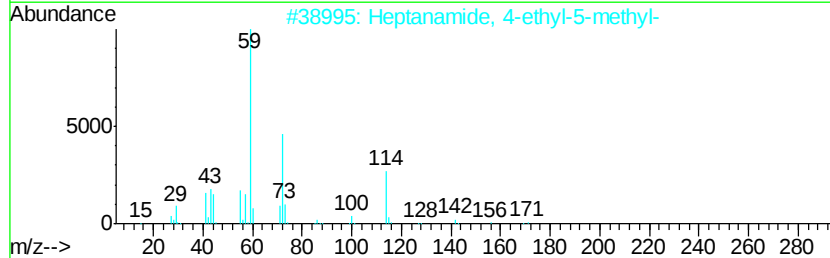
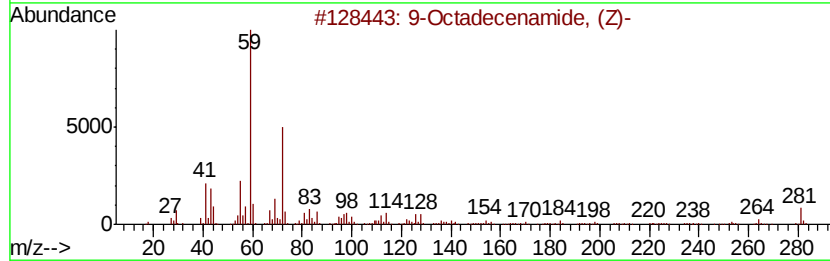
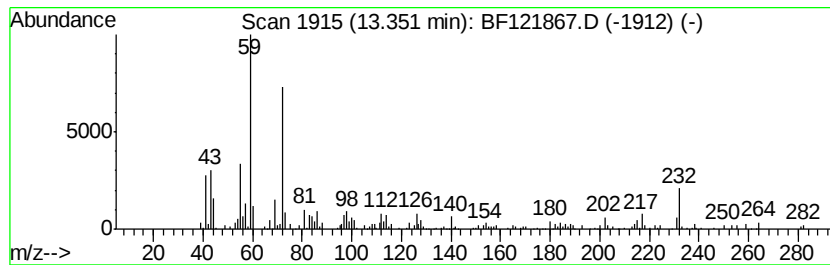
Instrument :
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ClientSampleId :
PAV-B18-092920-0-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 9-Octadecenamide, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.35	2.52 ng	140773	Chrysene-d12	13.92		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	90
2		Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1	64
3		Undecanamide, 11-bromo-	263	C11H22BrNO	005875-26-3	53
4		Dodecanamide	199	C12H25NO	001120-16-7	53
5		Nonanamide	157	C9H19NO	001120-07-6	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF100720\
Data File : BF121867.D
Acq On : 7 Oct 2020 11:47
Operator : JU/CG
Sample : L4190-11 2X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PAV-B18-092920-0-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF091020.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
5H-Indeno[1,2-b]p...	11.52	2.3	ng	150631	4	11.29	1296640	20.0
Phenanthrene, 1-m...	11.82	3.5	ng	227988	4	11.29	1296640	20.0
4H-Cyclopenta[def...	11.90	2.4	ng	156045	4	11.29	1296640	20.0
Octadecanoic acid	12.61	5.7	ng	317147	5	13.92	1116370	20.0
11H-Benzo[b]fluorene	13.05	3.6	ng	202108	5	13.92	1116370	20.0
9-Octadecenamide,...	13.35	2.5	ng	140773	5	13.92	1116370	20.0