

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042420\
 Data File : BF120198.D
 Acq On : 24 Apr 2020 13:23
 Operator : MA/SJ
 Sample : L2120-01
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-02-042220

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-BF040820.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	5.263	537	540	552	rBV	1943465	1910213	76.47%	6.821%
2	6.310	714	718	730	rBV	2161237	2013612	80.61%	7.190%
3	6.498	747	750	755	rBV	52640	50000	2.00%	0.179%
4	6.663	774	778	781	rBV	1147054	981179	39.28%	3.503%
5	6.816	800	804	808	rBV	2046925	1676632	67.12%	5.987%
6	7.228	869	874	877	rBV	1594979	1331395	53.30%	4.754%
7	7.945	992	996	999	rBV	1564134	1298745	51.99%	4.637%
8	8.657	1111	1117	1121	rVB	37963	38500	1.54%	0.137%
9	8.757	1131	1134	1138	rVB	45698	42031	1.68%	0.150%
10	9.028	1175	1180	1183	rBV	2870183	2497913	100.00%	8.919%
11	9.110	1190	1194	1199	rBV3	26071	40022	1.60%	0.143%
12	9.292	1221	1225	1228	rBV2	28363	38391	1.54%	0.137%
13	9.357	1233	1236	1239	rBV	57362	56052	2.24%	0.200%
14	9.422	1243	1247	1251	rVB	363142	287896	11.53%	1.028%
15	9.557	1267	1270	1276	rVB	68874	58489	2.34%	0.209%
16	9.698	1289	1294	1297	rBV	1623615	1372266	54.94%	4.900%
17	9.728	1297	1299	1304	rVB	62012	60321	2.41%	0.215%
18	9.904	1326	1329	1332	rVB	59096	46144	1.85%	0.165%
19	10.016	1343	1348	1351	rBV3	33761	39305	1.57%	0.140%
20	10.104	1359	1363	1365	rBV4	21396	28512	1.14%	0.102%
21	10.245	1384	1387	1393	rVB	110537	100891	4.04%	0.360%
22	10.404	1411	1414	1419	rVB4	26671	30512	1.22%	0.109%
23	10.486	1424	1428	1432	rBV	1607473	1359574	54.43%	4.855%
24	10.616	1446	1450	1454	rVB4	48311	58879	2.36%	0.210%
25	10.792	1476	1480	1483	rBV3	37804	46749	1.87%	0.167%
26	11.080	1527	1529	1534	rVB2	48795	55464	2.22%	0.198%
27	11.180	1542	1546	1548	rBV	1467091	1249055	50.00%	4.460%
28	11.204	1548	1550	1554	rVV	675190	493368	19.75%	1.762%
29	11.257	1554	1559	1562	rVB	157185	142702	5.71%	0.510%
30	11.292	1562	1565	1568	rBV	38179	41993	1.68%	0.150%
31	11.416	1583	1586	1593	rBV	42501	50618	2.03%	0.181%
32	11.510	1600	1602	1611	rVB2	42393	47186	1.89%	0.168%
33	11.586	1611	1615	1620	rVB2	64499	72316	2.90%	0.258%
34	11.680	1628	1631	1634	rBV	129129	121478	4.86%	0.434%

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Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040820.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.722	1634	1638	1642	rVV2	336979	432766	17.33%	1.545%
36	11.757	1642	1644	1646	rVV	79291	75171	3.01%	0.268%
37	11.792	1646	1650	1652	rVV	223778	243275	9.74%	0.869%
38	11.816	1652	1654	1659	rVB	114376	102702	4.11%	0.367%
39	11.974	1678	1681	1684	rBV	61961	62700	2.51%	0.224%
40	12.004	1684	1686	1689	rVB3	42017	40063	1.60%	0.143%
41	12.157	1710	1712	1714	rBV	50171	42804	1.71%	0.153%
42	12.180	1714	1716	1719	rVB2	41503	32505	1.30%	0.116%
43	12.233	1721	1725	1727	rBV	118821	114000	4.56%	0.407%
44	12.269	1727	1731	1733	rVV	167347	186062	7.45%	0.664%
45	12.292	1733	1735	1737	rVV	189413	156730	6.27%	0.560%
46	12.316	1737	1739	1743	rVB2	56762	67336	2.70%	0.240%
47	12.392	1748	1752	1756	rBV	849773	705750	28.25%	2.520%
48	12.439	1756	1760	1766	rVB6	87055	130771	5.24%	0.467%
49	12.504	1766	1771	1775	rBV4	119781	169166	6.77%	0.604%
50	12.616	1785	1790	1793	rBV	642510	694780	27.81%	2.481%
51	12.680	1798	1801	1804	rVB2	51487	57484	2.30%	0.205%
52	12.739	1808	1811	1812	rBV	45161	35113	1.41%	0.125%
53	12.769	1812	1816	1819	rVV	2337466	1924742	77.05%	6.873%
54	12.839	1823	1828	1831	rBV4	78881	116556	4.67%	0.416%
55	12.927	1839	1843	1844	rBV4	53164	63801	2.55%	0.228%
56	12.945	1844	1846	1850	rVB	147463	145276	5.82%	0.519%
57	13.016	1850	1858	1861	rBV2	145547	190573	7.63%	0.680%
58	13.051	1861	1864	1867	rBV2	101919	98345	3.94%	0.351%
59	13.145	1877	1880	1882	rVV2	65779	52151	2.09%	0.186%
60	13.174	1882	1885	1888	rVB	72702	62890	2.52%	0.225%
61	13.257	1896	1899	1904	rVB4	85245	79061	3.17%	0.282%
62	13.457	1931	1933	1937	rVB	42837	46332	1.85%	0.165%
63	13.563	1947	1951	1954	rBV2	66436	93154	3.73%	0.333%
64	13.616	1958	1960	1962	rVV	43218	40064	1.60%	0.143%
65	13.674	1966	1970	1978	rVV3	167356	270548	10.83%	0.966%
66	13.816	1989	1994	1997	rBV2	1045140	1190639	47.67%	4.251%
67	13.839	1997	1998	2001	rVB	244757	195471	7.83%	0.698%
68	13.921	2010	2012	2014	rVB	59452	45027	1.80%	0.161%
69	14.204	2058	2060	2063	rVB	87335	69247	2.77%	0.247%
70	14.239	2063	2066	2068	rBV	46230	49470	1.98%	0.177%
71	14.298	2074	2076	2079	rBV3	92526	92858	3.72%	0.332%

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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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72	14.827	2162	2166	2169	rBV	276995	392568	15.72%	1.402%
73	14.904	2177	2179	2182	rBV	66283	83705	3.35%	0.299%
74	15.104	2210	2213	2219	rVB	158462	199434	7.98%	0.712%
75	15.163	2219	2223	2228	rBV	251176	293372	11.74%	1.048%
76	15.221	2228	2233	2236	rBV	698935	815406	32.64%	2.912%
77	15.662	2304	2308	2311	rBV3	63833	100642	4.03%	0.359%
78	16.562	2457	2461	2470	rVB2	113827	207405	8.30%	0.741%

Sum of corrected areas: 28006318

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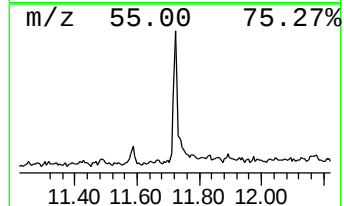
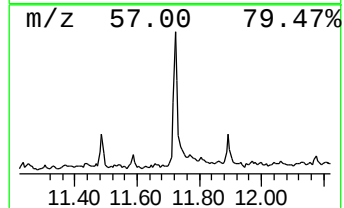
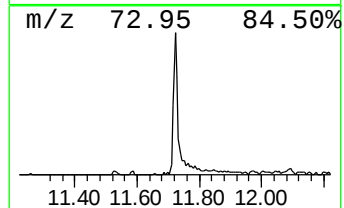
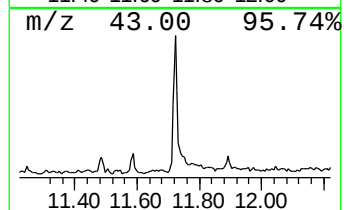
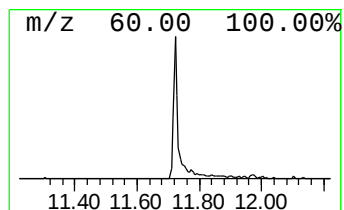
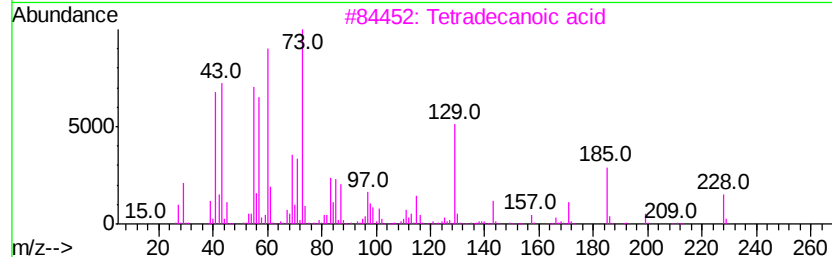
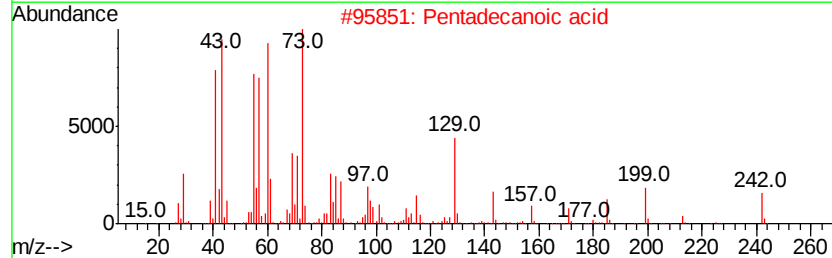
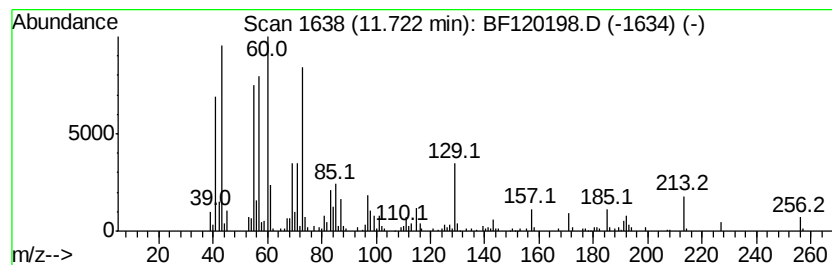
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.72	6.93 ng	432766	Phenanthrene-d10	11.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	86
4		Tridecanoic acid	214	C13H26O2	000638-53-9	72
5		Dodecanoic acid	200	C12H24O2	000143-07-7	64



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Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
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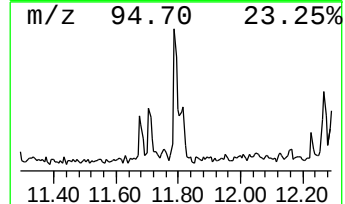
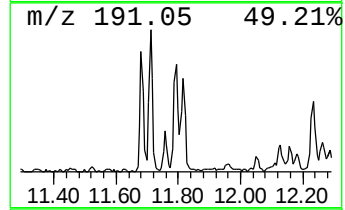
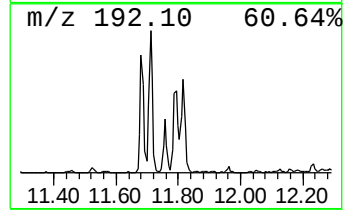
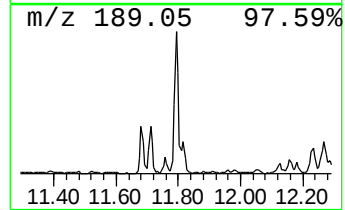
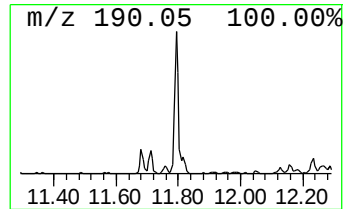
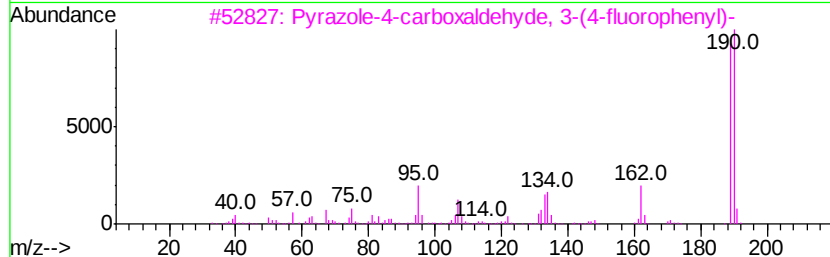
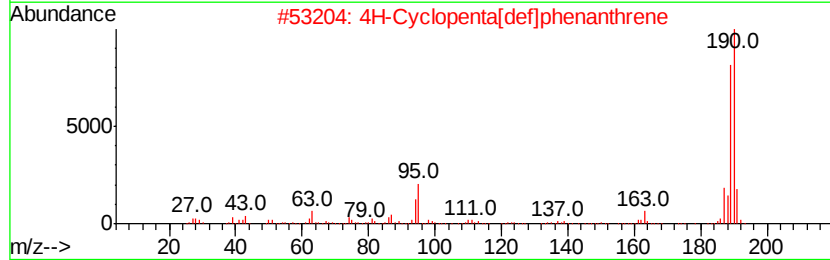
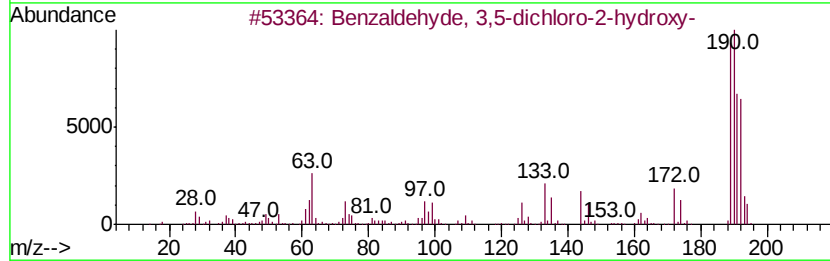
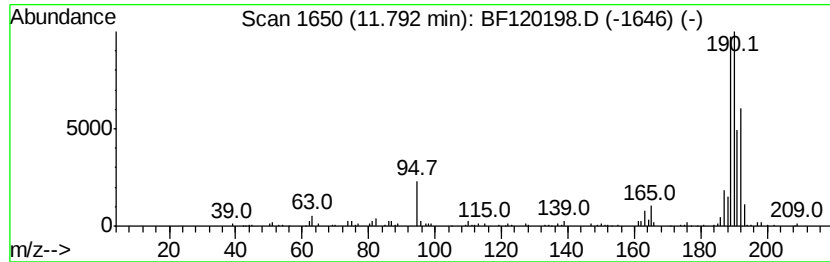
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Benzaldehyde, 3,5-dichloro-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	3.90 ng	243275	Phenanthrene-d10	11.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	72
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
3		Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	49
4		Naphtho[1,2-b]norborene	192	C15H12	1000210-14-8	38
5		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	25



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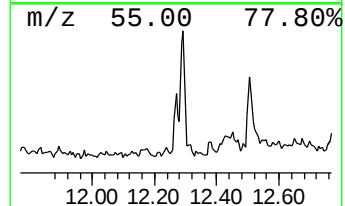
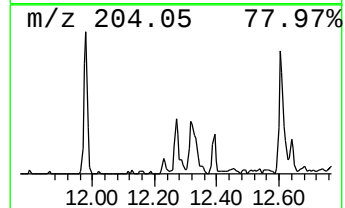
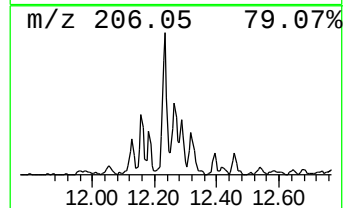
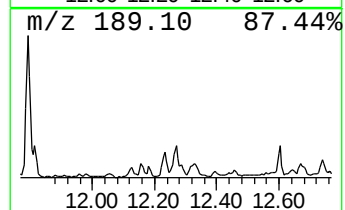
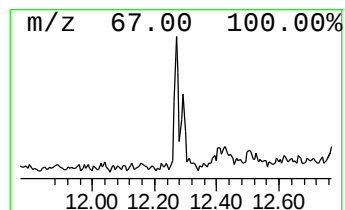
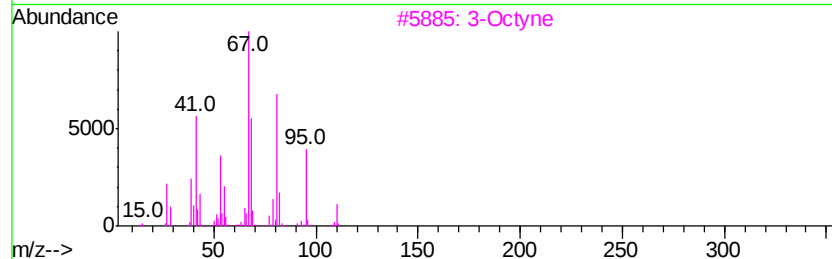
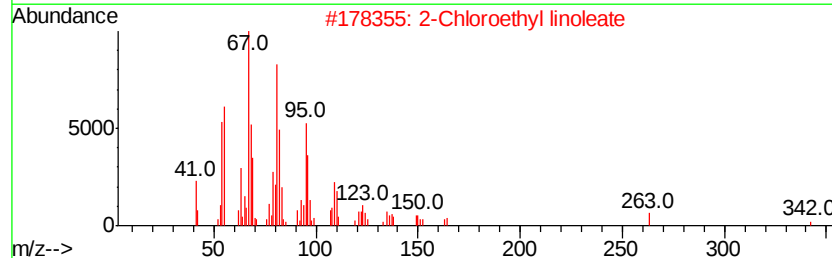
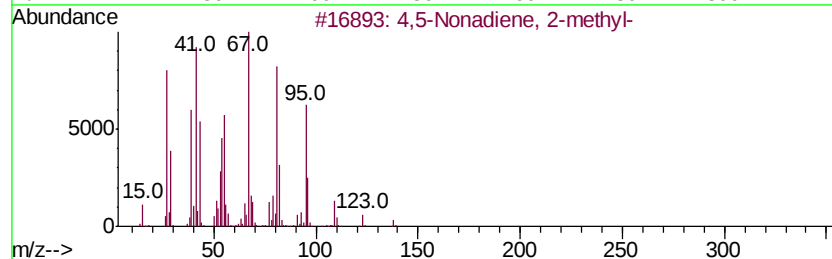
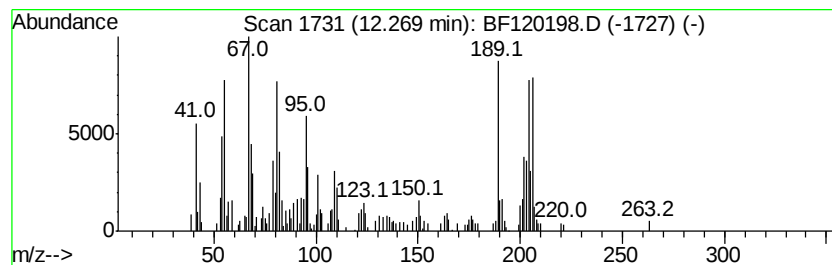
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown12.27 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.27	2.98 ng	186062	Phenanthrene-d10	11.18

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,5-Nonadiene, 2-methyl-			138	C10H18	055956-32-6	38
2	2-Chloroethyl linoleate			342	C20H35ClO2	025525-76-2	38
3	3-Octyne			110	C8H14	015232-76-5	30
4	4-(Trifluoromethoxy)benzoic acid			206	C8H5F3O3	000330-12-1	25
5	4H-Thiazolo[5,4-b]indole, 7-fluo...			206	C10H7FN2S	1000337-12-3	25



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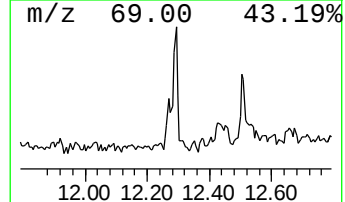
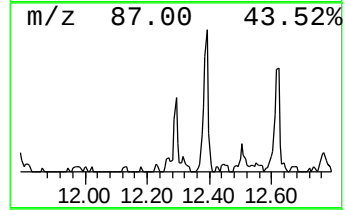
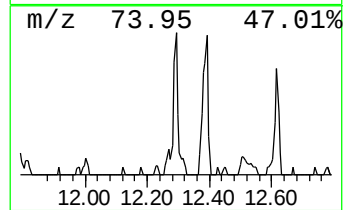
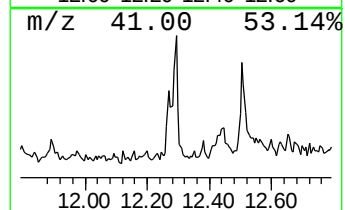
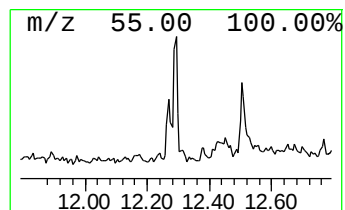
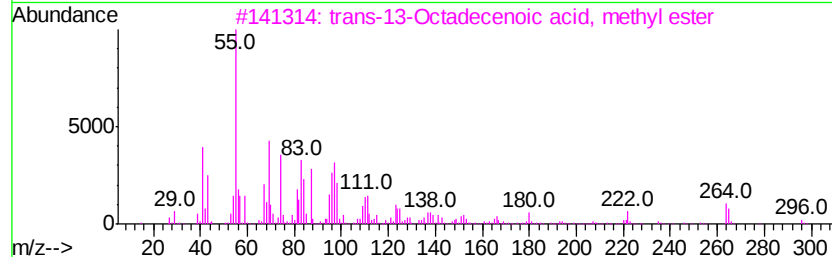
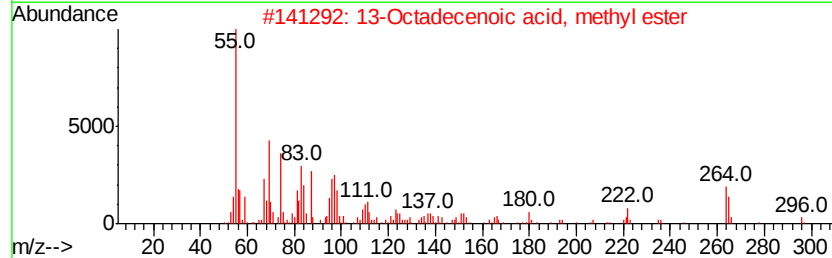
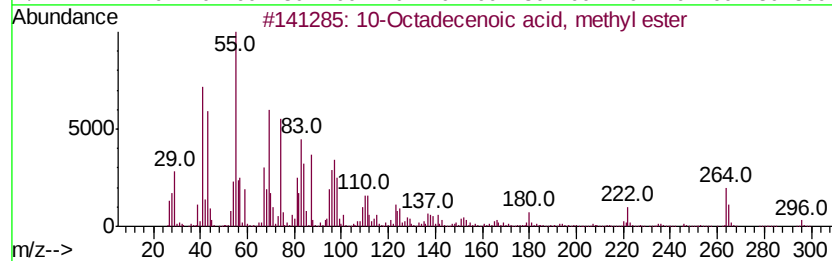
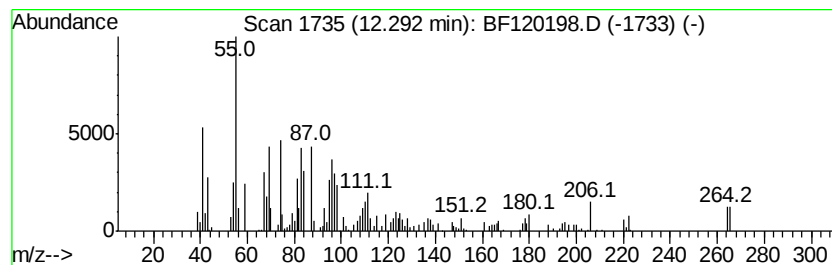
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 10-Octadecenoic acid, methyl ester Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	2.51 ng	156730	Phenanthrene-d10	11.18

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	87
2		13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	86
3		trans-13-Octadecenoic acid, methyl ester	296	C19H36O2	1000333-61-3	68
4		9-Octadecenoic acid (Z)-, methyl ester	296	C19H36O2	000112-62-9	64
5		14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042420\
Data File : BF120198.D
Acq On : 24 Apr 2020 13:23
Operator : MA/SJ
Sample : L2120-01
Misc :
ALS Vial : 35 Sample Multiplier: 1

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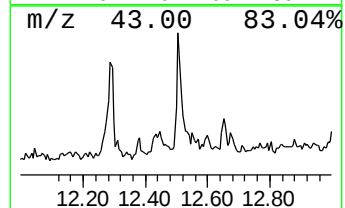
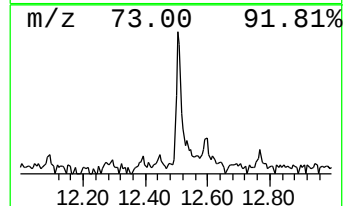
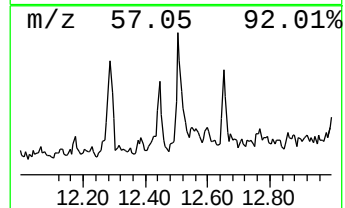
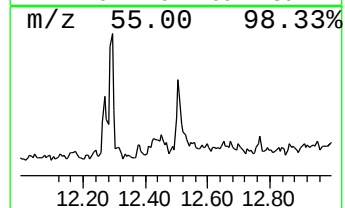
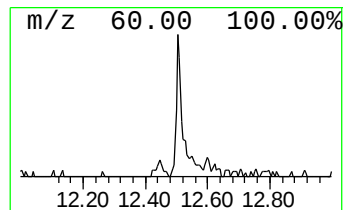
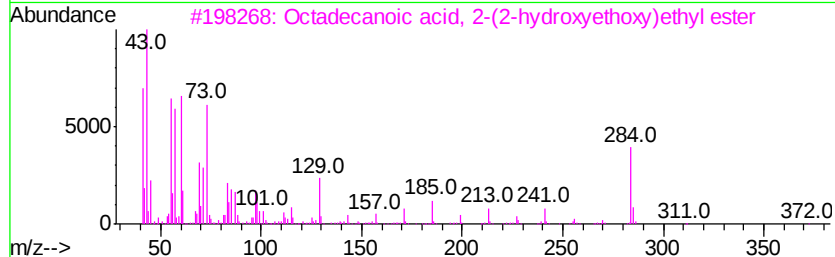
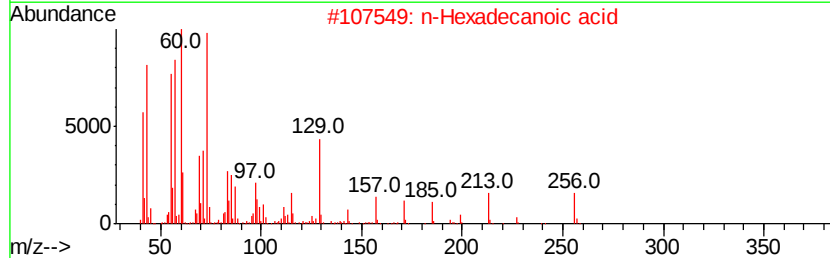
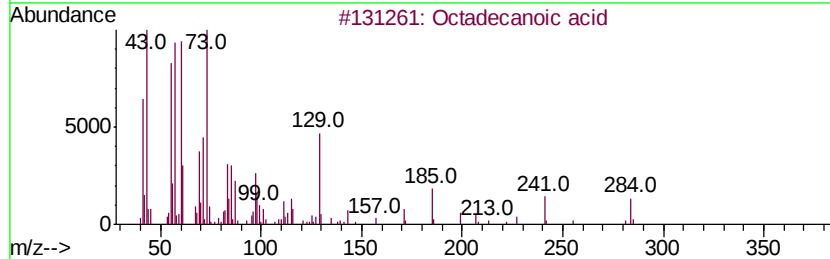
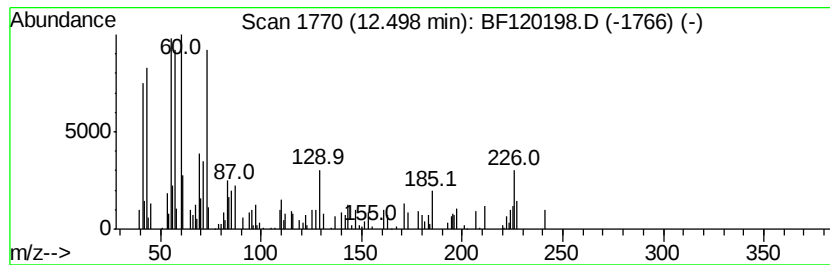
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.50	2.84 ng	169166	Chrysene-d12	13.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	97
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
3		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	90
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5		Decanoic acid, silver(1+) salt	278	C10H19AgO2	013126-67-5	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042420\
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ClientSampleId :
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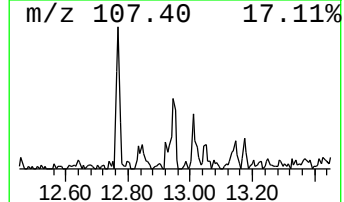
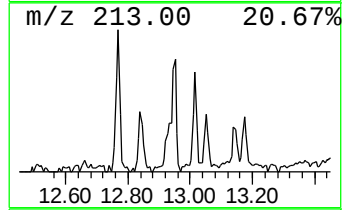
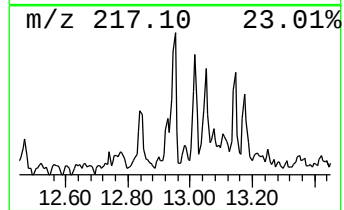
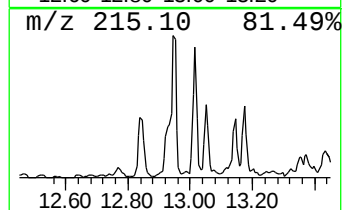
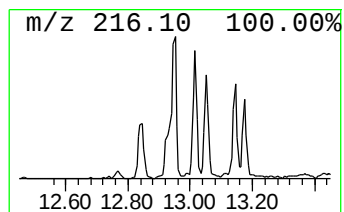
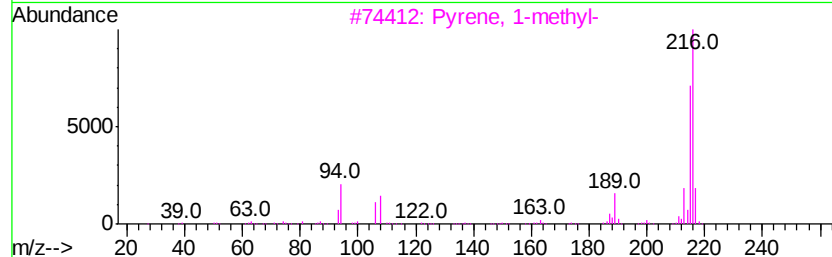
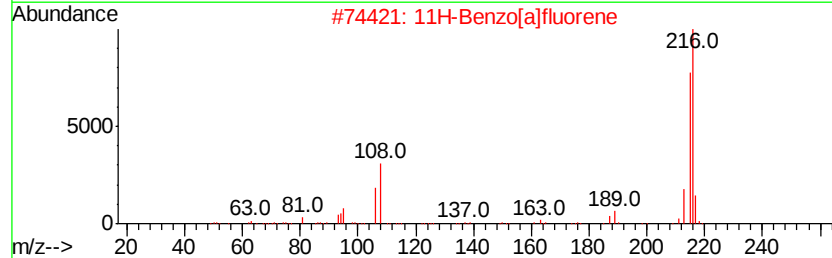
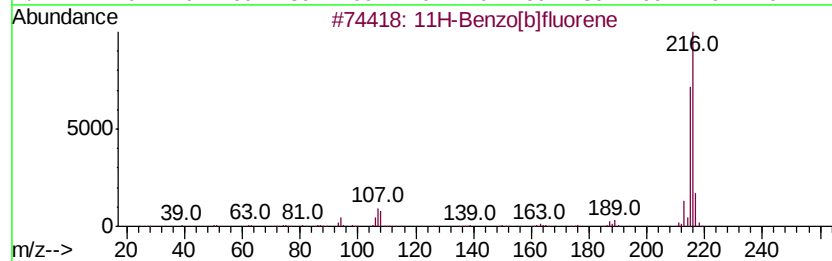
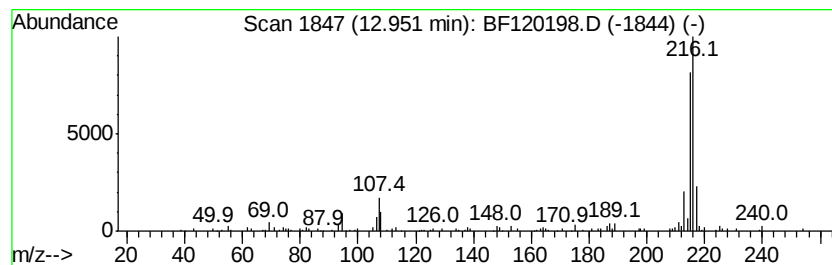
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 11H-Benzo[b]fluorene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.95	2.44 ng	145276	Chrysene-d12	13.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
2		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	83
3		Pvrene, 1-methyl-	216	C17H12	002381-21-7	74
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	72
5		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042420\
Data File : BF120198.D
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Operator : MA/SJ
Sample : L2120-01
Misc :
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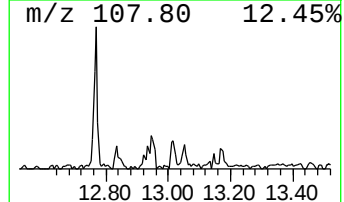
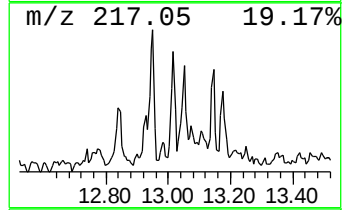
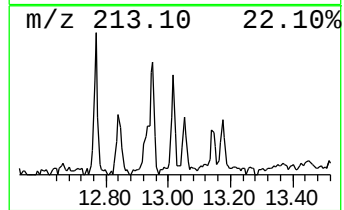
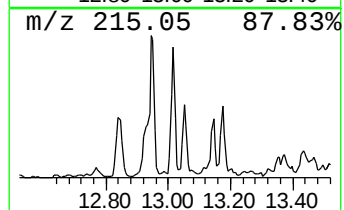
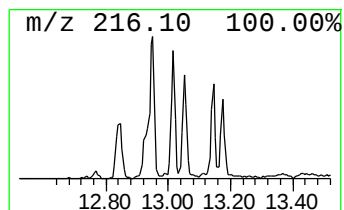
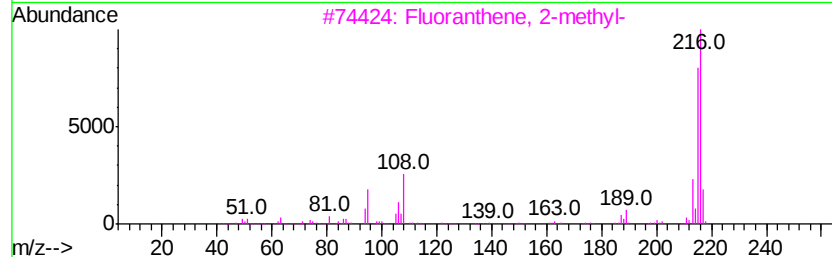
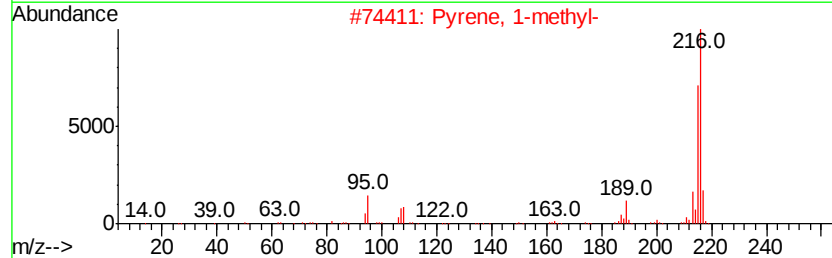
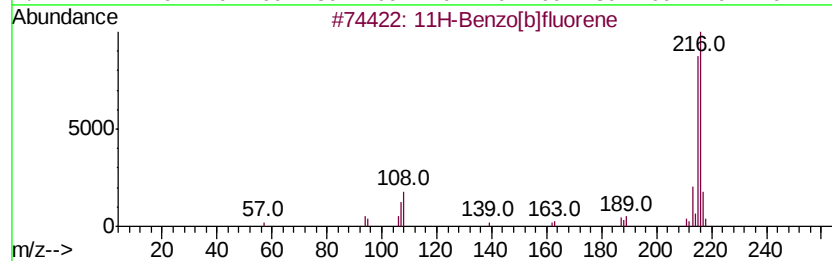
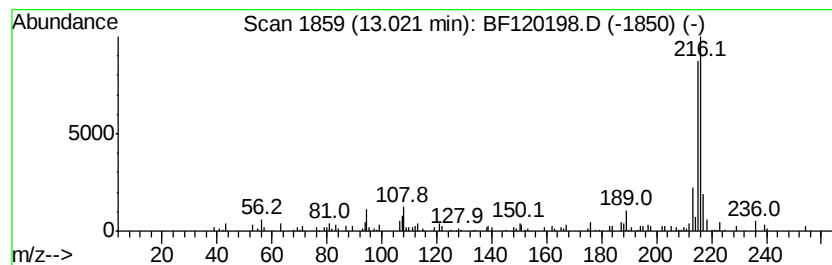
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Pyrene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.02	3.20 ng	190573	Chrysene-d12	13.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[blfluorene	216	C17H12	000243-17-4	95
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
4	11H-Benzo[alfluorene	216	C17H12	000238-84-6	83
5	1,3,5-Triazine-2,4,6-triamine, N...	216	C10H12N6	046731-79-7	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042420\
Data File : BF120198.D
Acq On : 24 Apr 2020 13:23
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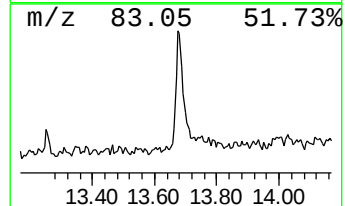
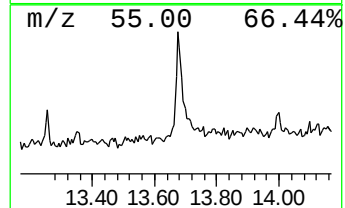
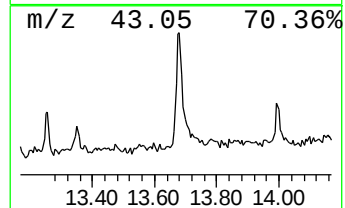
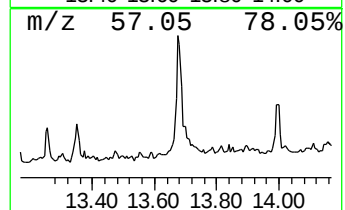
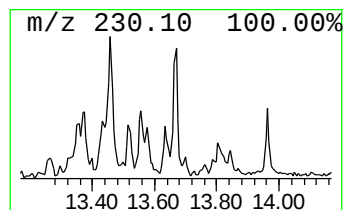
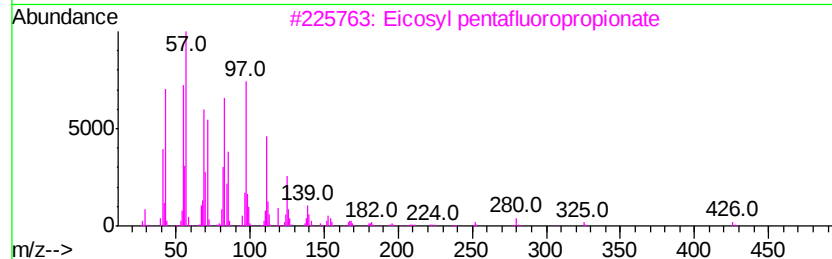
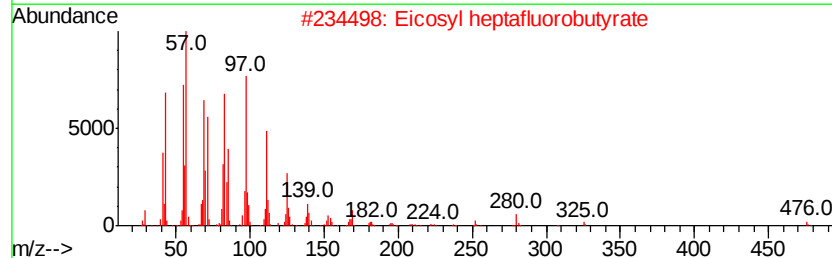
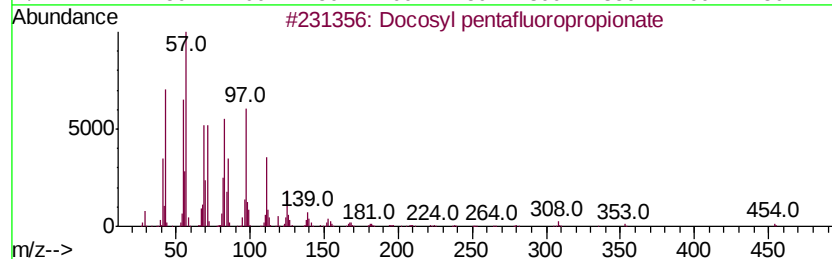
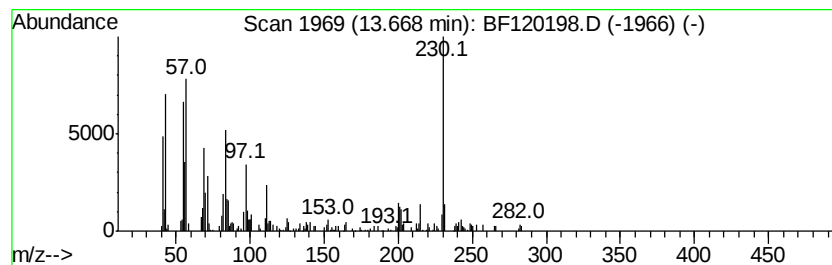
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Docosyl pentafluoropropionate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	4.54 ng	270548	Chrysene-d12	13.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	87
2		Eicosyl heptafluorobutyrate	494	C24H41F7O2	1000351-83-5	86
3		Eicosyl pentafluoropropionate	444	C23H41F5O2	1000351-80-8	86
4		17-Pentatriacontene	491	C35H70	006971-40-0	83
5		n-Heptadecanol-1	256	C17H36O	001454-85-9	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042420\
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Sample : L2120-01
Misc :
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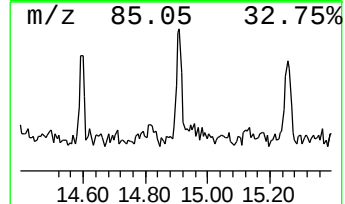
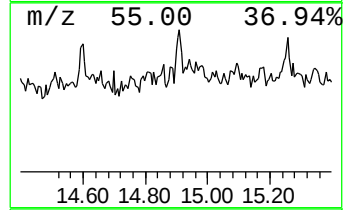
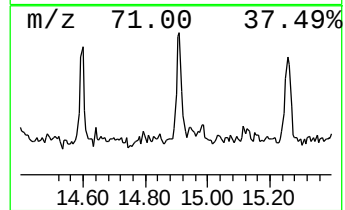
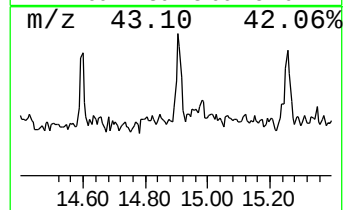
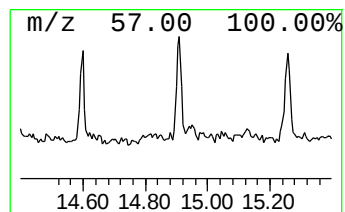
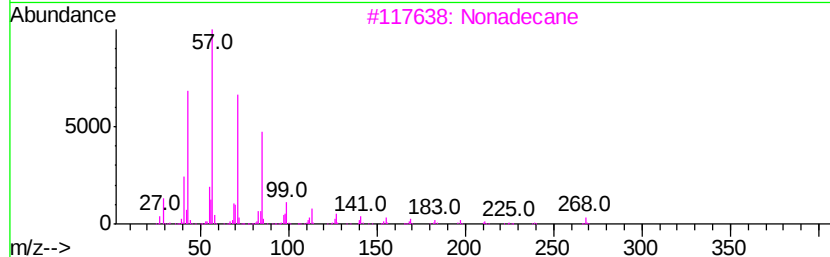
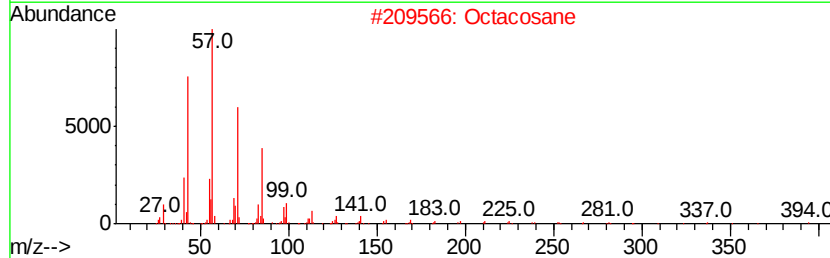
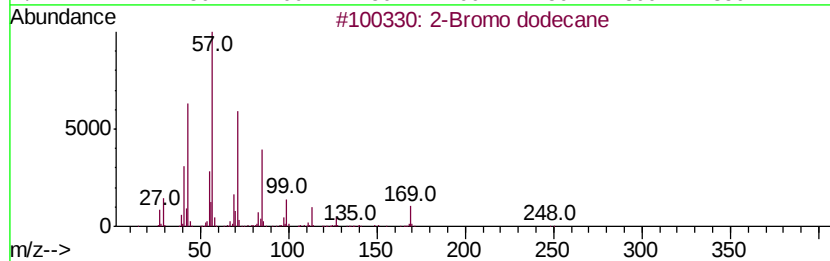
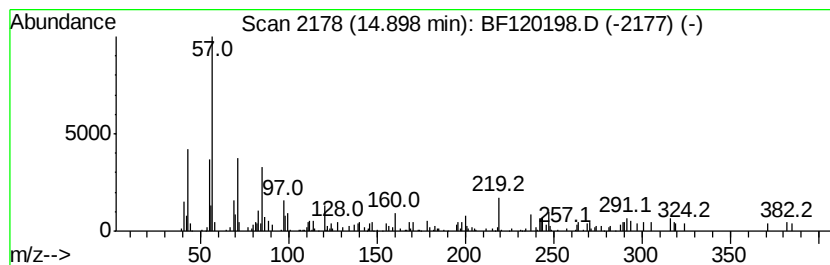
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2-Bromo dodecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.90	2.05 ng	83705	Perylene-d12	15.22	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromo dodecane	248	C12H25Br	013187-99-0	87
2	Octacosane	394	C28H58	000630-02-4	83
3	Nonadecane	268	C19H40	000629-92-5	80
4	Tetracosane	338	C24H50	000646-31-1	80
5	Tetradecane	198	C14H30	000629-59-4	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042420\
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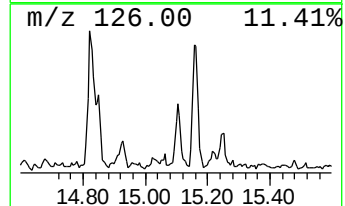
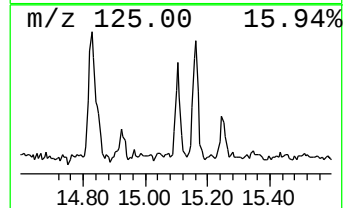
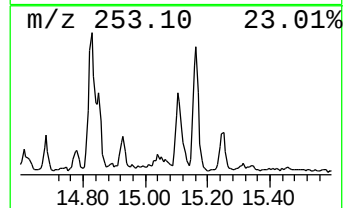
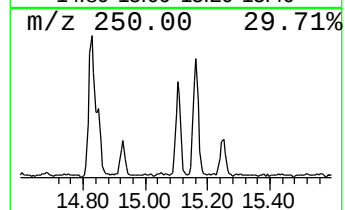
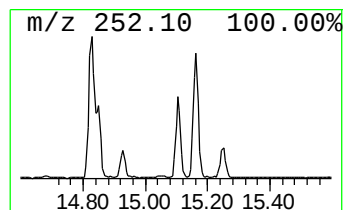
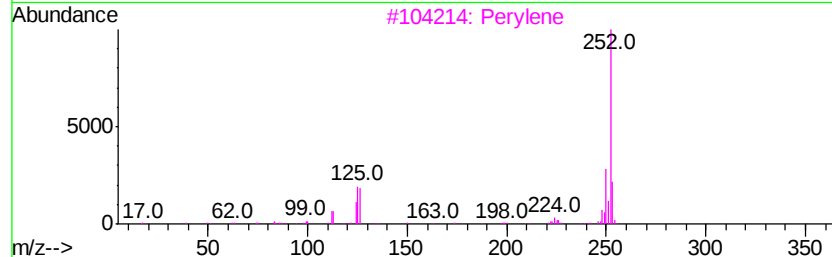
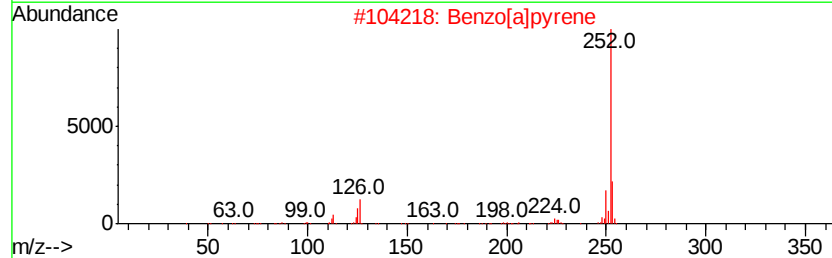
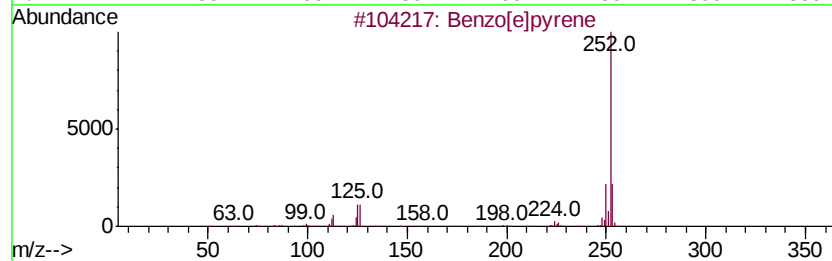
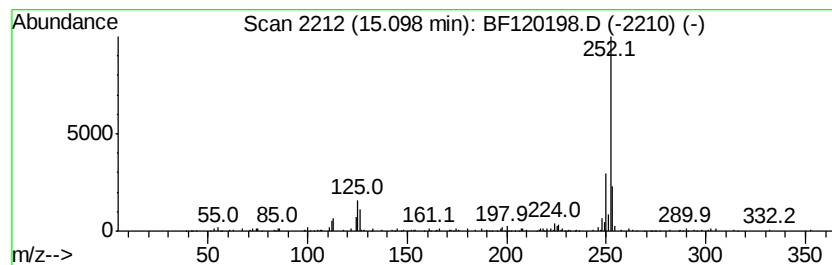
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.10	4.89 ng	199434	Perylene-d12	15.22

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[a]pyrene	252	C20H12	000050-32-8	96
3			Perylene	252	C20H12	000198-55-0	95
4			Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	93
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042420\
Data File : BF120198.D
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Operator : MA/SJ
Sample : L2120-01
Misc :
ALS Vial : 35 Sample Multiplier: 1

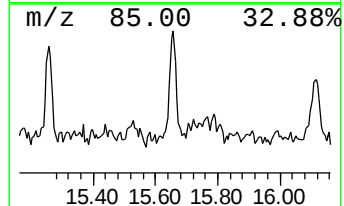
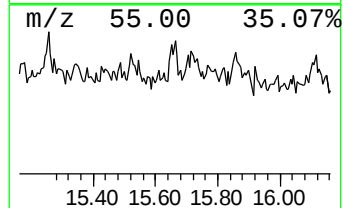
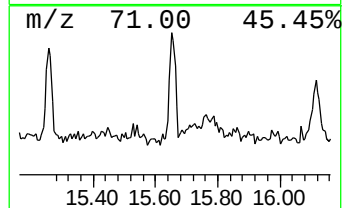
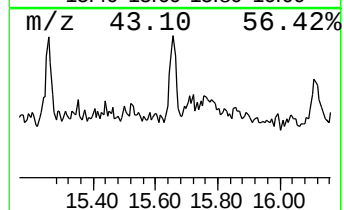
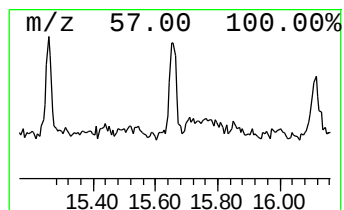
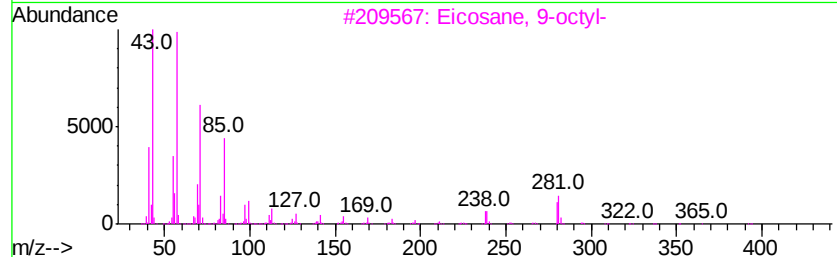
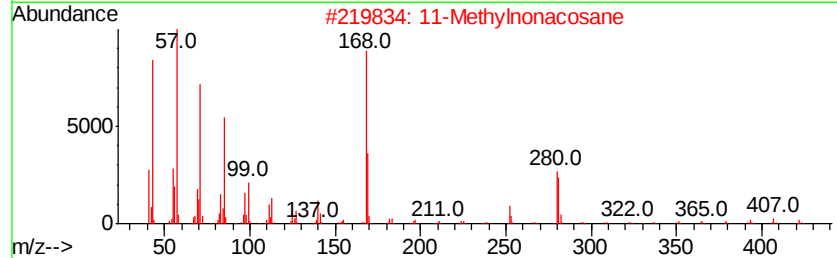
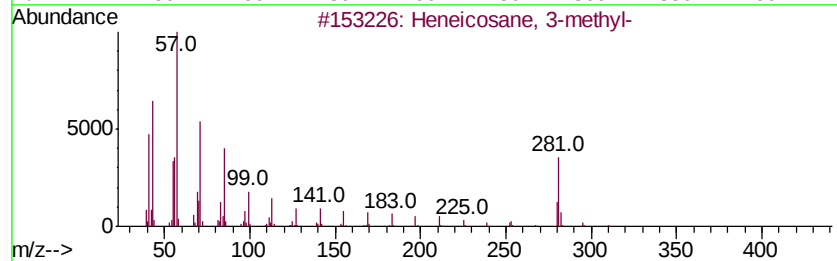
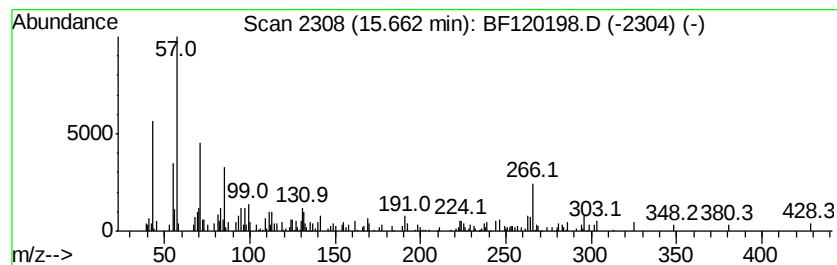
Instrument :
BNA_F
ClientSampleId :
OR-02-042220

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown15.66 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	2.47 ng	100642	Perylene-d12	15.22
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Heneicosane, 3-methyl-	310 C22H46	006418-47-9 35
2		11-Methylnonacosane	422 C30H62	007371-99-5 35
3		Eicosane, 9-octyl-	394 C28H58	013475-77-9 35
4		Dodecane, 1-iodo-	296 C12H25I	004292-19-7 32
5		Hexadecanoic acid, 2-(octadecylo...	553 C36H72O3	029899-13-6 32



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042420\
 Data File : BF120198.D
 Acq On : 24 Apr 2020 13:23
 Operator : MA/SJ
 Sample : L2120-01
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-02-042220

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040820.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
n-Hexadecanoic acid	11.72	6.9 ng		432766	4	11.18	1249060	20.0
Benzaldehyde, 3,5...	11.79	3.9 ng		243275	4	11.18	1249060	20.0
unknown12.27	12.27	3.0 ng		186062	4	11.18	1249060	20.0
10-Octadecenoic a...	12.29	2.5 ng		156730	4	11.18	1249060	20.0
Octadecanoic acid	12.50	2.8 ng		169166	5	13.82	1190640	20.0
11H-Benzo[b]fluorene	12.95	2.4 ng		145276	5	13.82	1190640	20.0
Pyrene, 1-methyl-	13.02	3.2 ng		190573	5	13.82	1190640	20.0
Docosyl pentafluo...	13.67	4.5 ng		270548	5	13.82	1190640	20.0
2-Bromo dodecane	14.90	2.0 ng		83705	6	15.22	815406	20.0
Benzo[e]pyrene	15.10	4.9 ng		199434	6	15.22	815406	20.0
unknown15.66	15.66	2.5 ng		100642	6	15.22	815406	20.0