

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF010919\
 Data File : BF111963.D
 Acq On : 9 Jan 2019 21:31
 Operator : JU/SJ
 Sample : K1044-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CPSB-11(20-25)

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-BF010719.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	5.099	507	512	520	rVB	127645	153620	1.03%	0.173%
2	5.516	579	583	592	rVB	2603207	2340978	15.69%	2.642%
3	6.169	691	694	697	rBV	268243	221981	1.49%	0.251%
4	6.528	750	755	760	rBV	2715319	2667385	17.88%	3.011%
5	6.657	773	777	781	rVB	2640763	2471184	16.56%	2.789%
6	6.881	812	815	823	rVB	877439	838022	5.62%	0.946%
7	7.040	838	842	846	rBV	2252118	1953962	13.10%	2.206%
8	7.440	902	910	915	rBV	1614758	1713699	11.49%	1.934%
9	7.481	915	917	921	rVB	278963	213118	1.43%	0.241%
10	7.904	987	989	996	rVB5	159286	289937	1.94%	0.327%
11	8.169	1028	1034	1036	rVV2	1110228	1653239	11.08%	1.866%
12	8.187	1036	1037	1040	rVV	268156	221244	1.48%	0.250%
13	8.228	1040	1044	1047	rVB2	217861	259583	1.74%	0.293%
14	8.463	1078	1084	1088	rBV4	244443	329825	2.21%	0.372%
15	8.516	1090	1093	1096	rVB3	183829	164078	1.10%	0.185%
16	8.551	1096	1099	1101	rBV3	195555	223032	1.50%	0.252%
17	8.587	1101	1105	1107	rBV2	250734	207615	1.39%	0.234%
18	8.669	1116	1119	1122	rBV2	383085	361363	2.42%	0.408%
19	8.757	1131	1134	1138	rBV	1093257	887653	5.95%	1.002%
20	8.992	1169	1174	1176	rBV2	339377	371848	2.49%	0.420%
21	9.122	1192	1196	1200	rVB3	195081	326854	2.19%	0.369%
22	9.186	1205	1207	1209	rVV2	279825	233828	1.57%	0.264%
23	9.239	1212	1216	1219	rBV	3057089	2567761	17.21%	2.898%
24	9.322	1226	1230	1235	rBV	1167608	1113560	7.46%	1.257%
25	9.392	1240	1242	1245	rVB	255833	225065	1.51%	0.254%
26	9.563	1267	1271	1272	rBV4	137845	178549	1.20%	0.202%
27	9.581	1272	1274	1276	rVV	315491	280672	1.88%	0.317%
28	9.604	1276	1278	1280	rVV2	214793	189865	1.27%	0.214%
29	9.633	1280	1283	1287	rVV2	569846	841813	5.64%	0.950%
30	9.669	1287	1289	1291	rVV2	259255	230489	1.55%	0.260%
31	9.704	1291	1295	1297	rVB	268669	264125	1.77%	0.298%
32	9.798	1306	1311	1313	rBV5	96560	164857	1.11%	0.186%
33	9.851	1313	1320	1323	rVV	1115769	1088168	7.29%	1.228%
34	9.922	1329	1332	1335	rVV2	1089230	964298	6.46%	1.088%

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35	9.957	1335	1338	1340	rVB2	198132	175136	1.17%	0.198%
36	10.086	1357	1360	1362	rVV3	123810	171598	1.15%	0.194%
37	10.122	1362	1366	1371	rVB5	150135	303689	2.04%	0.343%
38	10.175	1371	1375	1379	rBV3	204141	283882	1.90%	0.320%
39	10.239	1383	1386	1389	rBV2	166984	217159	1.46%	0.245%
40	10.351	1403	1405	1408	rVB	1083179	849929	5.70%	0.959%
41	10.469	1422	1425	1431	rBV4	175074	228962	1.53%	0.258%
42	10.575	1438	1443	1448	rBV	360398	508332	3.41%	0.574%
43	10.710	1464	1466	1470	rVB	1588906	1306951	8.76%	1.475%
44	10.828	1482	1486	1491	rBV	961045	1186753	7.96%	1.340%
45	11.275	1559	1562	1565	rBV	865959	684471	4.59%	0.773%
46	11.304	1565	1567	1572	rVB	422618	439587	2.95%	0.496%
47	11.410	1582	1585	1587	rBV	1015328	865503	5.80%	0.977%
48	11.433	1587	1589	1596	rVV	537171	566524	3.80%	0.639%
49	11.698	1631	1634	1637	rBV	634834	592717	3.97%	0.669%
50	11.804	1648	1652	1655	rBV	6996087	6187646	41.48%	6.984%
51	11.945	1672	1676	1681	rBV2	820652	887014	5.95%	1.001%
52	12.110	1701	1704	1708	rBV	565687	511872	3.43%	0.578%
53	12.157	1709	1712	1717	rVB3	267288	298808	2.00%	0.337%
54	12.269	1728	1731	1735	rBV2	437773	435604	2.92%	0.492%
55	12.357	1743	1746	1749	rVB	364989	301344	2.02%	0.340%
56	12.504	1762	1771	1772	rBV	10367361	14746484	98.85%	16.645%
57	12.527	1772	1775	1780	rVV	12177406	14918315	100.00%	16.839%
58	12.598	1783	1787	1790	rVV	3691457	3245840	21.76%	3.664%
59	12.669	1796	1799	1807	rVB	2582502	2914755	19.54%	3.290%
60	12.727	1807	1809	1815	rVB3	233652	243721	1.63%	0.275%
61	12.827	1824	1826	1829	rBV4	190310	180899	1.21%	0.204%
62	12.857	1829	1831	1839	rVB2	490185	722741	4.84%	0.816%
63	12.998	1852	1855	1858	rVB	2482867	2106018	14.12%	2.377%
64	13.139	1876	1879	1882	rBV	1053942	879991	5.90%	0.993%
65	13.227	1891	1894	1895	rBV	298470	294873	1.98%	0.333%
66	13.316	1907	1909	1912	rVB	368906	313730	2.10%	0.354%
67	13.569	1949	1952	1954	rBV	227445	186094	1.25%	0.210%
68	13.727	1976	1979	1985	rVB3	504171	553707	3.71%	0.625%
69	13.892	2004	2007	2013	rVB	305415	313028	2.10%	0.353%
70	13.986	2020	2023	2028	rBV	369753	367909	2.47%	0.415%
71	14.057	2031	2035	2037	rBV2	829432	846751	5.68%	0.956%

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72	14.210	2059	2061	2064	rVB	226348	181593	1.22%	0.205%
73	14.516	2111	2113	2119	rVB	222716	225589	1.51%	0.255%
74	14.827	2164	2166	2170	rVB	205403	174303	1.17%	0.197%
75	14.951	2182	2187	2196	rVB	559422	875864	5.87%	0.989%
76	15.174	2222	2225	2228	rVB	229536	244336	1.64%	0.276%
77	15.545	2284	2288	2298	rVB2	434016	841214	5.64%	0.950%

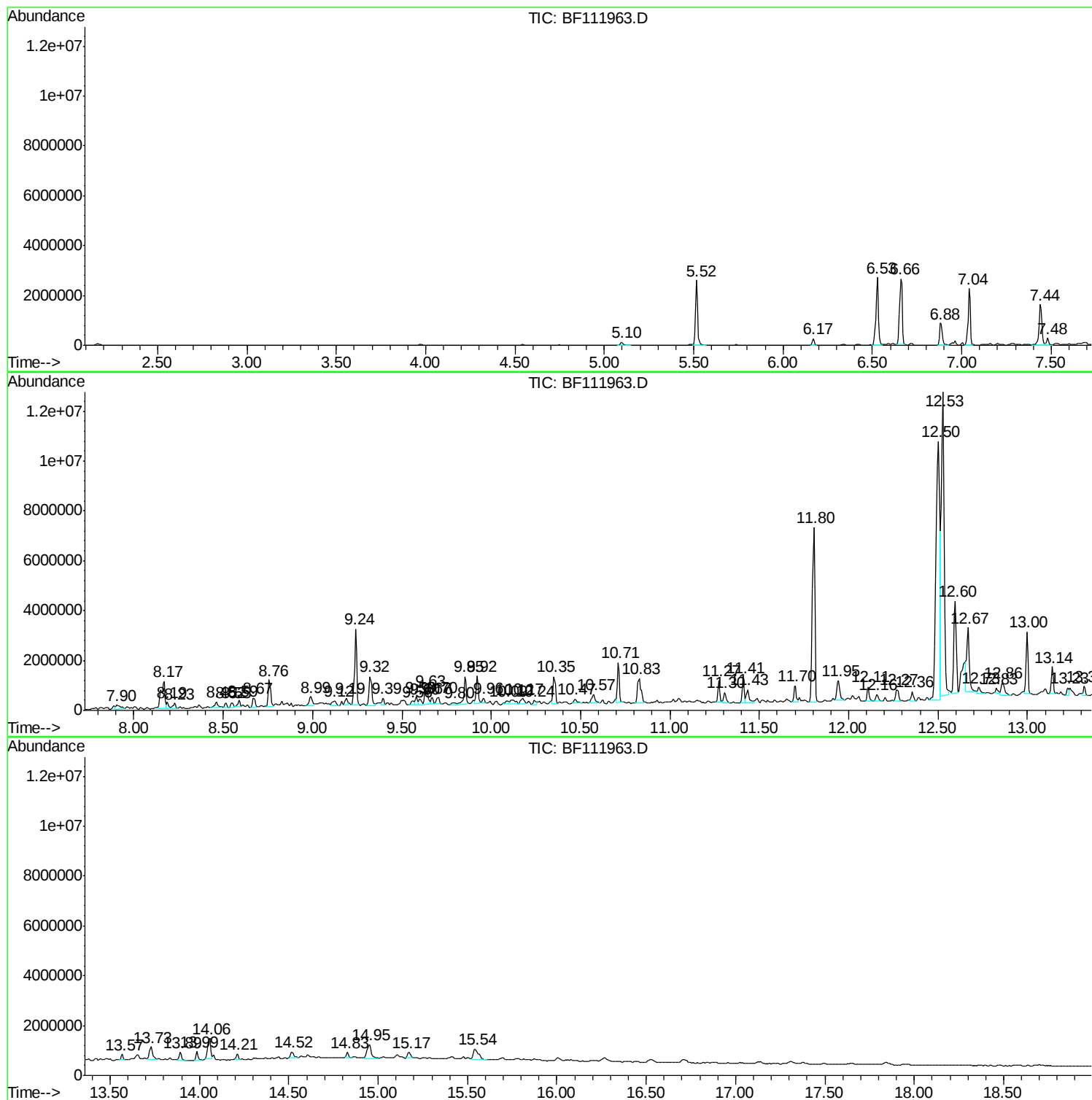
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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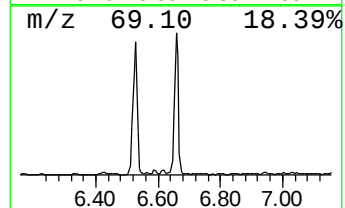
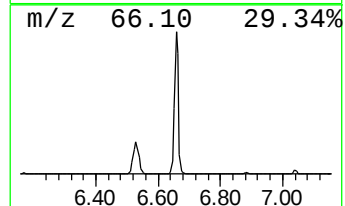
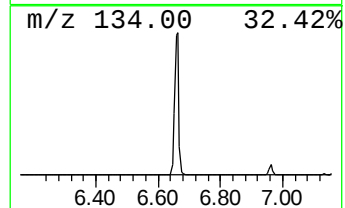
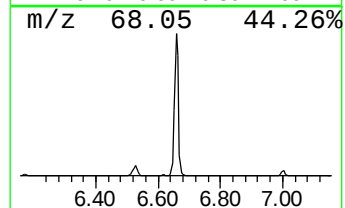
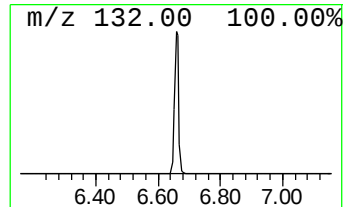
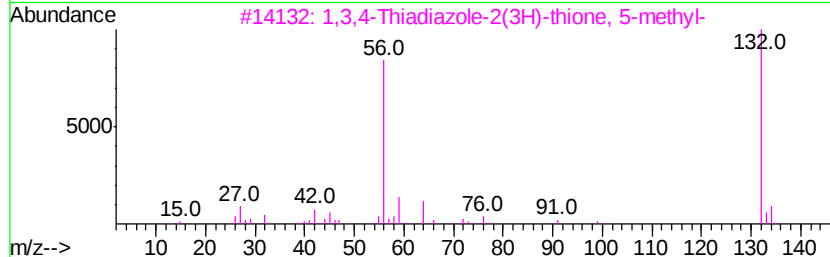
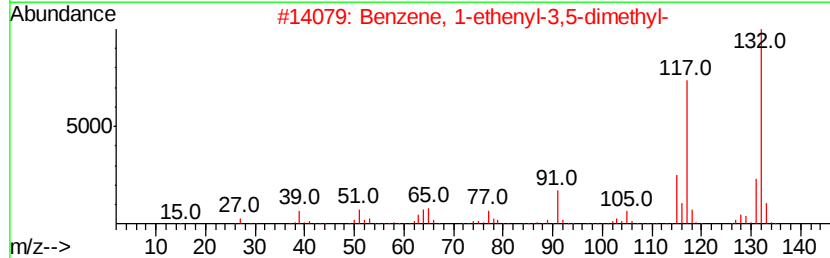
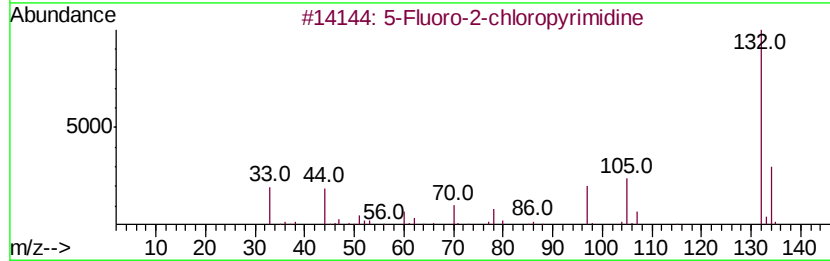
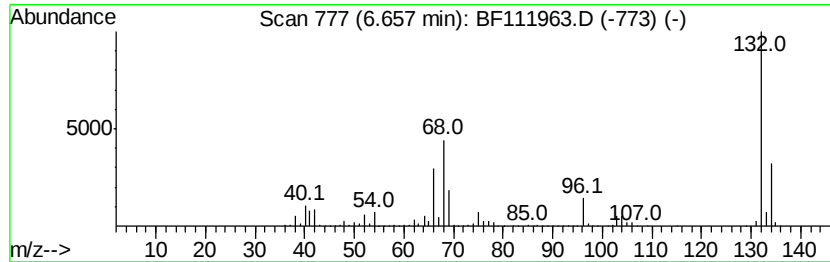
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.66 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	58.98 ng	2471180	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
3		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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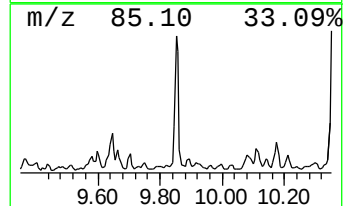
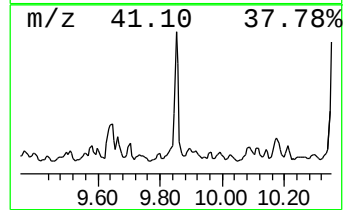
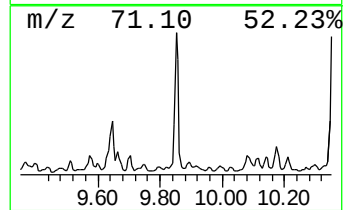
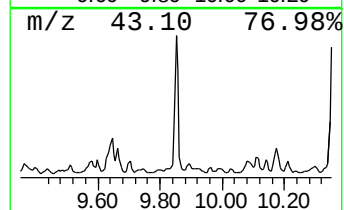
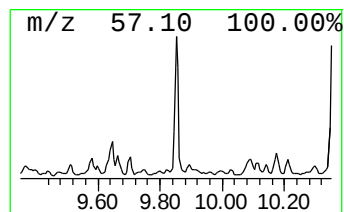
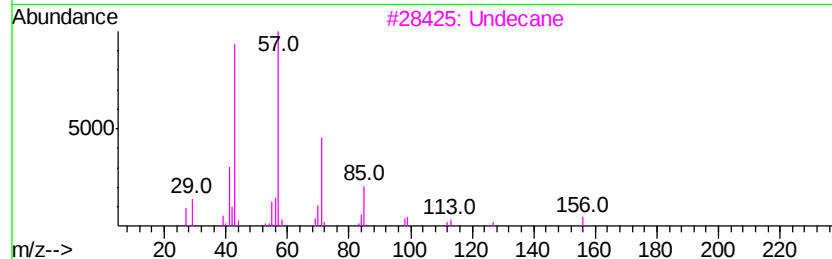
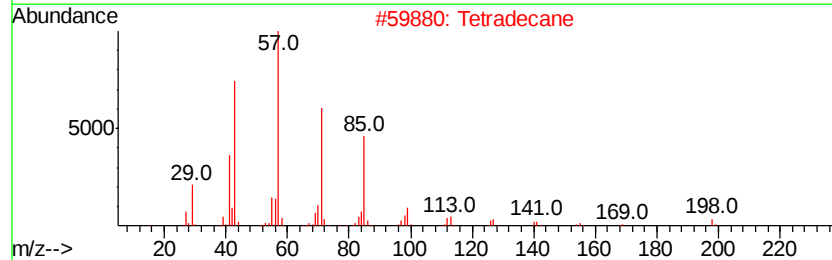
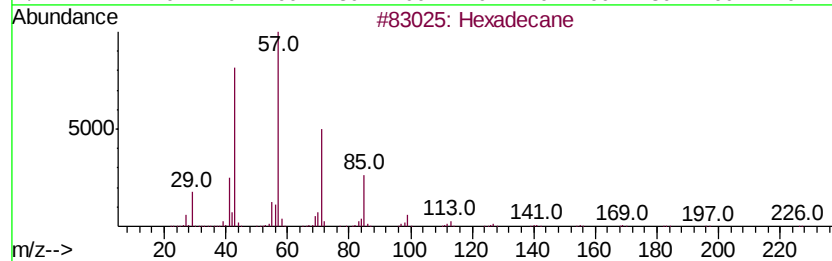
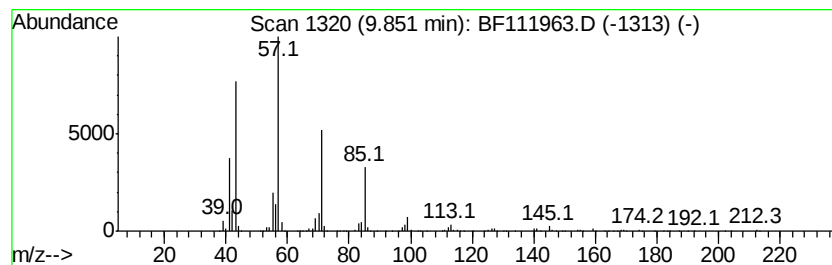
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Hexadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.85	22.57 ng	1088170	Acenaphthene-d10	9.92

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	91
2	Tetradecane		198	C14H30	000629-59-4	90
3	Undecane		156	C11H24	001120-21-4	86
4	Decane, 6-ethyl-2-methyl-		184	C13H28	062108-21-8	83
5	Tridecane		184	C13H28	000629-50-5	78



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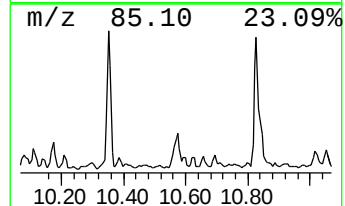
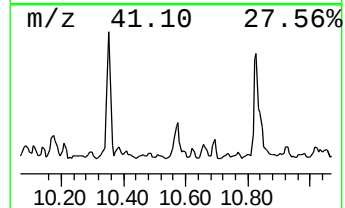
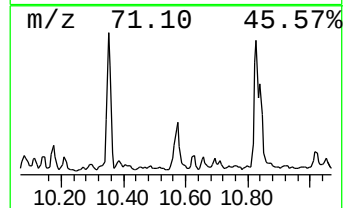
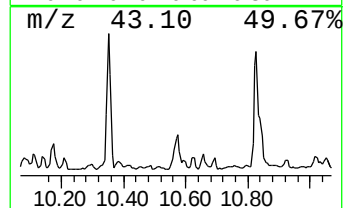
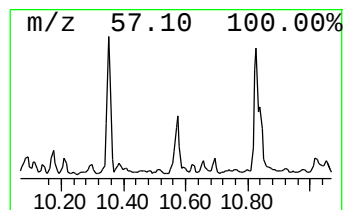
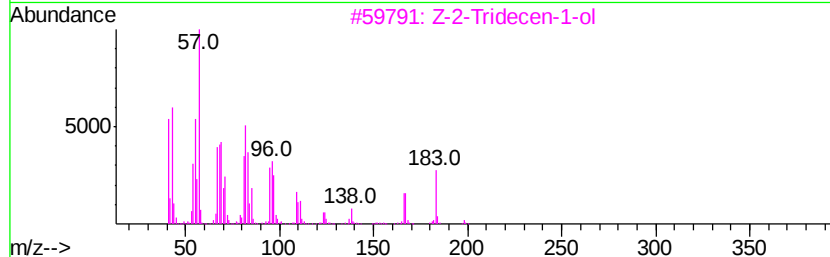
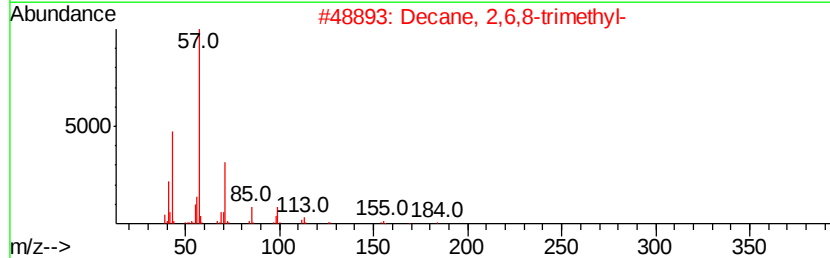
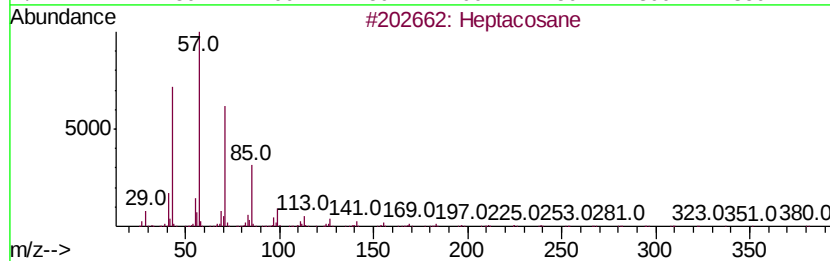
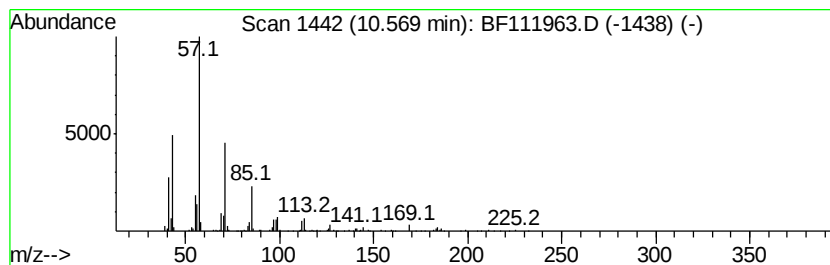
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Heptacosane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.57	10.54 ng	508332	Acenaphthene-d10	9.92

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	80
2			Decane, 2,6,8-trimethyl-	184	C13H28	062108-26-3	76
3			Z-2-Tridecen-1-ol	198	C13H26O	1000130-75-8	74
4			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	72
5			Hexadecane	226	C16H34	000544-76-3	64



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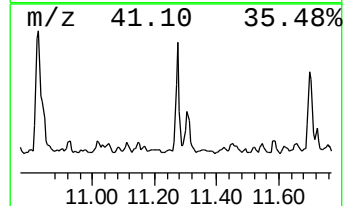
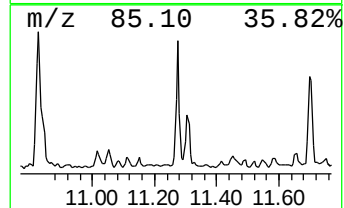
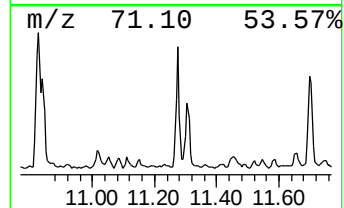
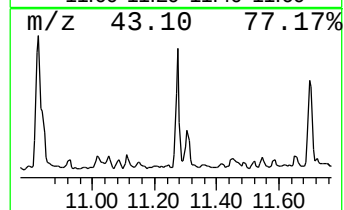
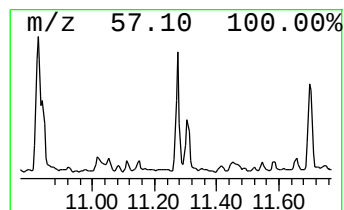
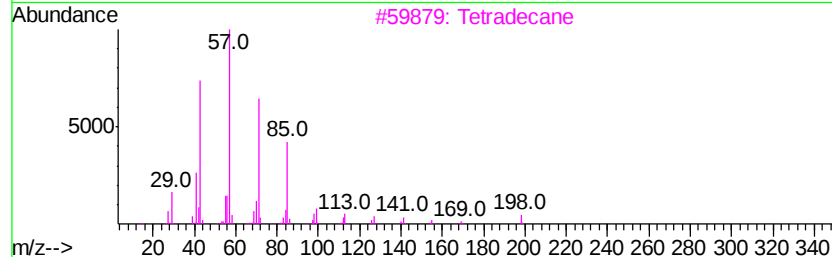
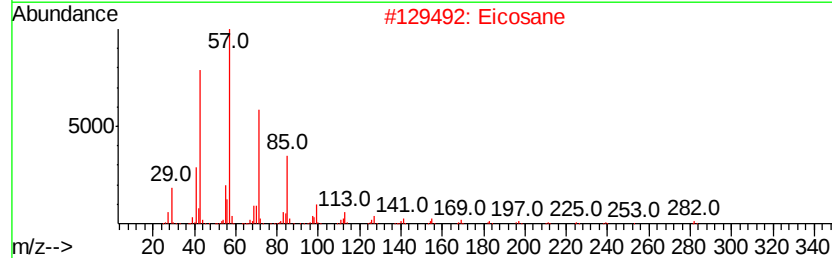
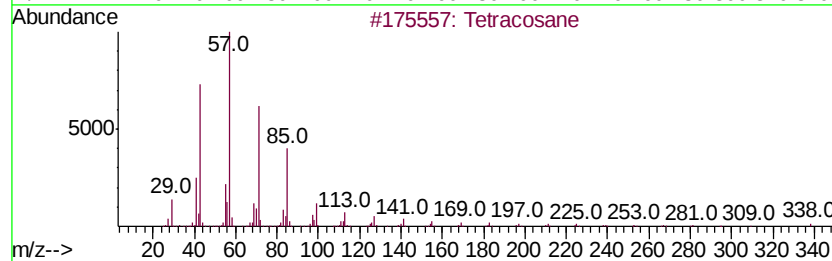
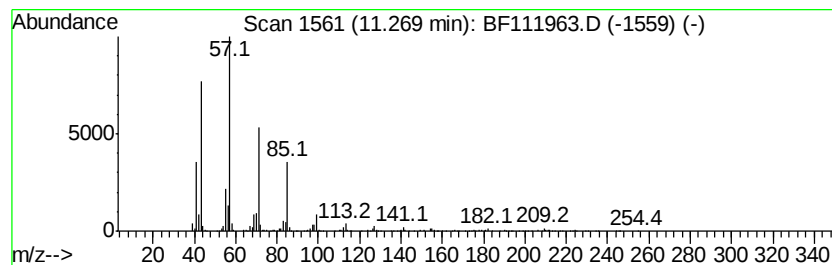
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetracosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.27	15.82 ng	684471	Phenanthrene-d10	11.41

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	90
2		Eicosane	282	C20H42	000112-95-8	90
3		Tetradecane	198	C14H30	000629-59-4	90
4		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90
5		Pentadecane	212	C15H32	000629-62-9	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010919\
Data File : BF111963.D
Acq On : 9 Jan 2019 21:31
Operator : JU/SJ
Sample : K1044-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CPSB-11(20-25)

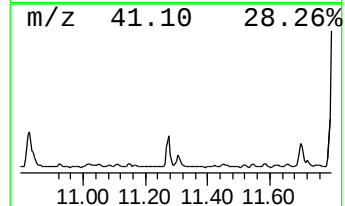
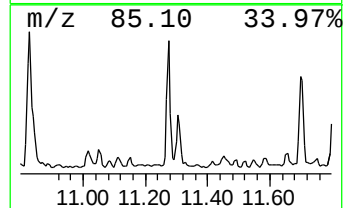
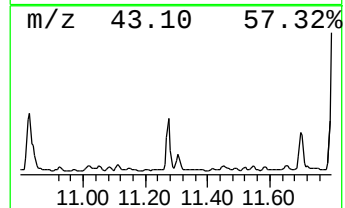
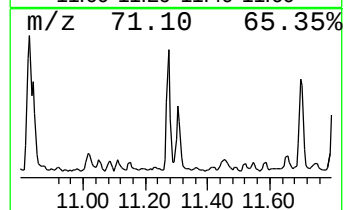
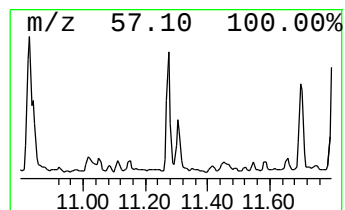
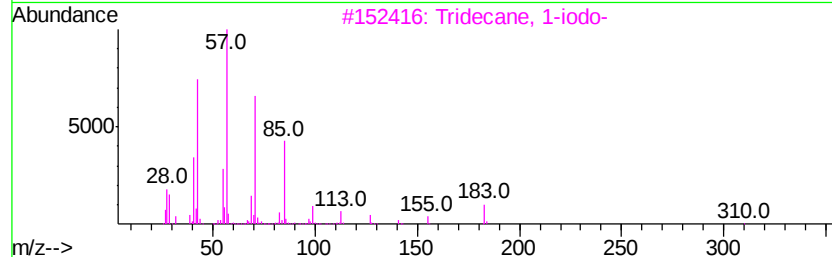
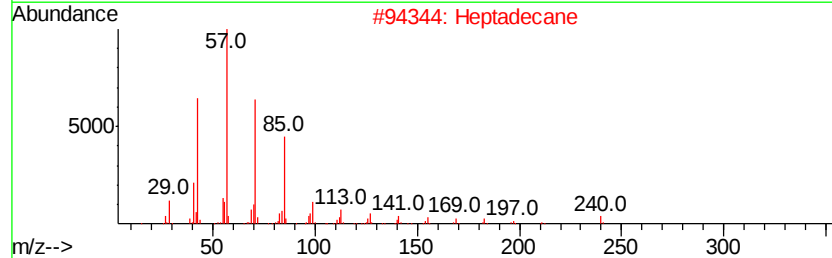
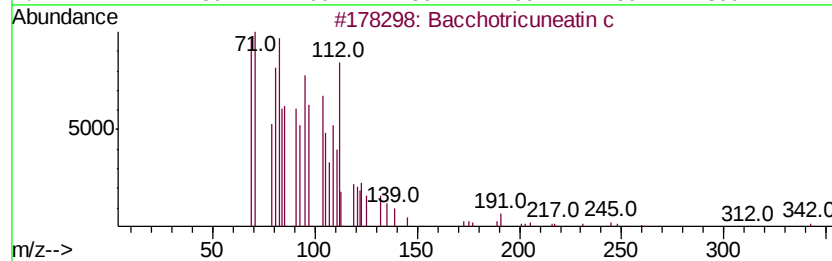
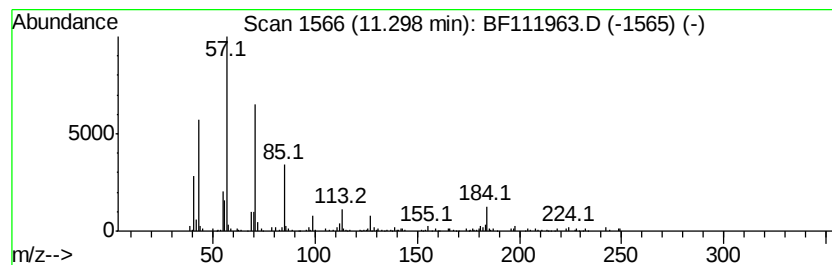
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Bacchotricuneatin c Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.30	10.16 ng	439587	Phenanthrene-d10	11.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bacchotricuneatin c	342	C20H22O5	066563-30-2	96
2		Heptadecane	240	C17H36	000629-78-7	72
3		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	64
4		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	64
5		3-Methyl-4-(methoxycarbonyl)hexa...	184	C9H12O4	1000104-10-8	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010919\
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Sample : K1044-04
Misc :
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ClientSampleId :
CPSB-11(20-25)

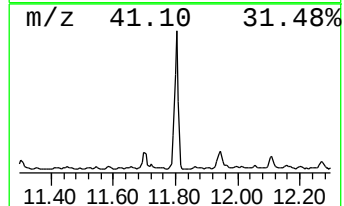
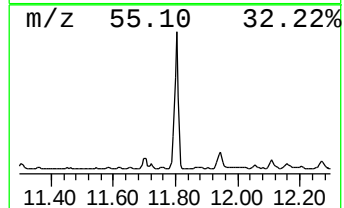
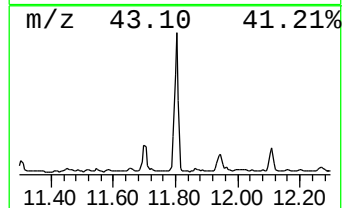
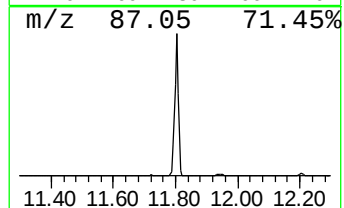
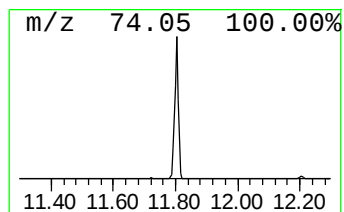
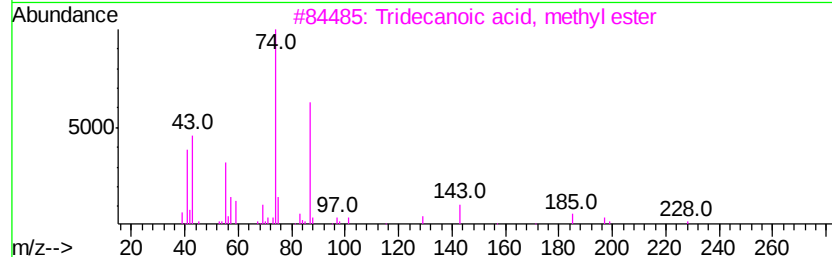
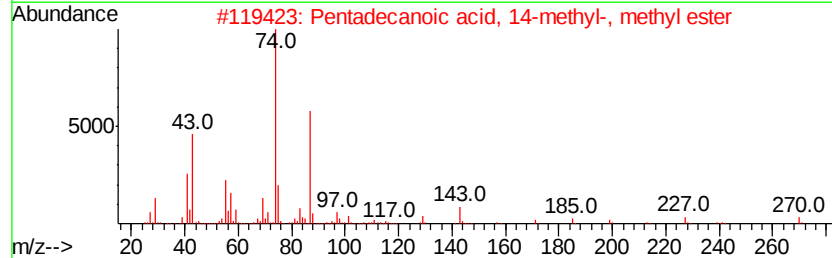
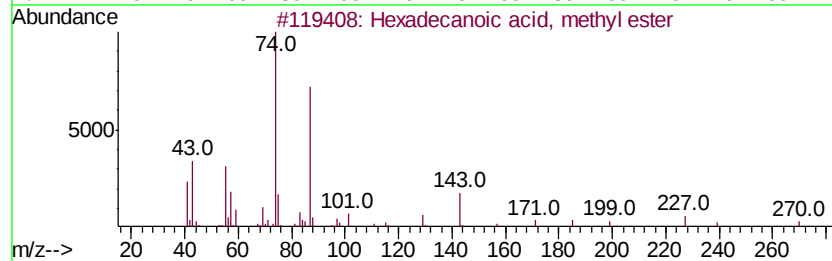
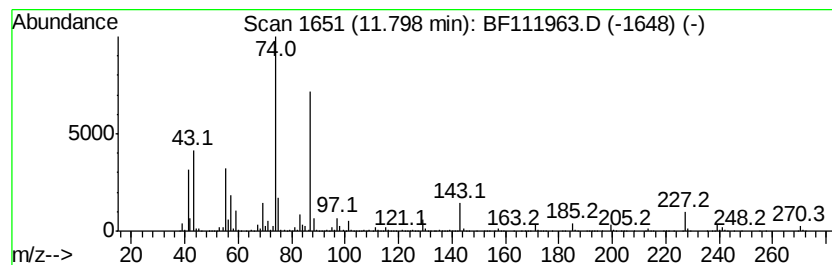
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.80	142.98 ng	6187650	Phenanthrene-d10	11.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	97
3		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	91
4		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	90
5		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010919\
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Misc :
ALS Vial : 17 Sample Multiplier: 1

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ClientSampleId :
CPSB-11(20-25)

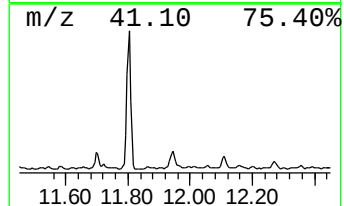
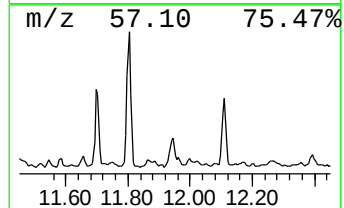
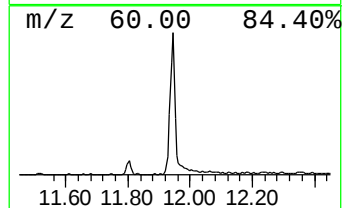
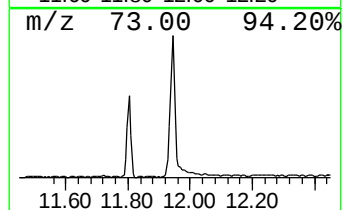
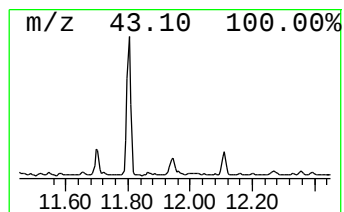
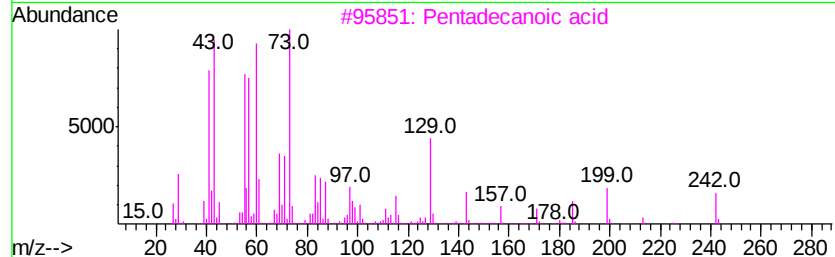
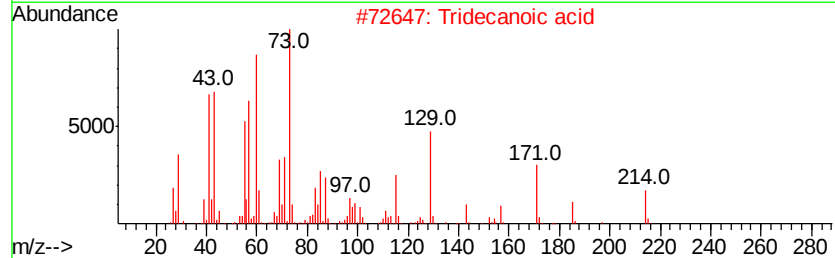
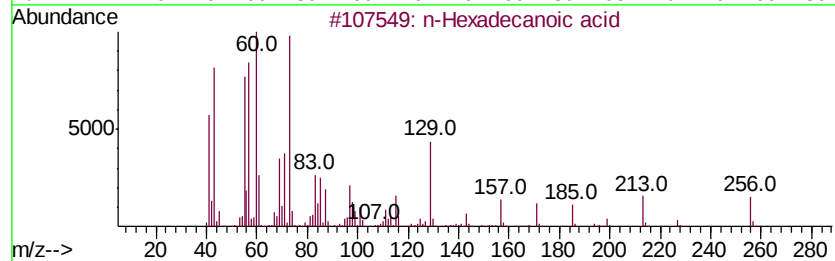
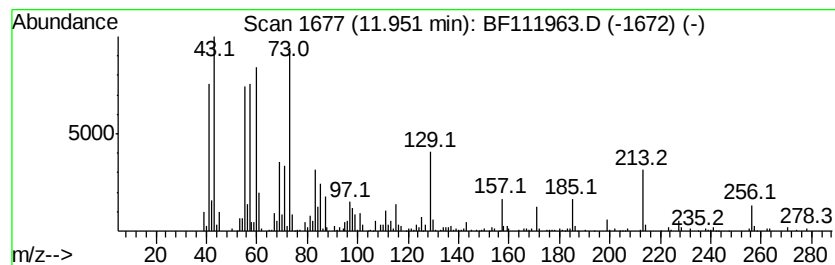
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.95	20.50 ng	887014	Phenanthrene-d10	11.41

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Tridecanoic acid	214	C13H26O2	000638-53-9	90
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	89
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	86
5			n-Decanoic acid	172	C10H20O2	000334-48-5	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010919\
Data File : BF111963.D
Acq On : 9 Jan 2019 21:31
Operator : JU/SJ
Sample : K1044-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CPSB-11(20-25)

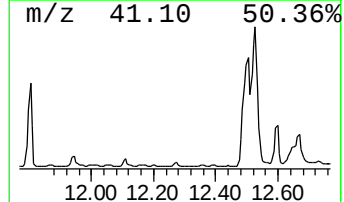
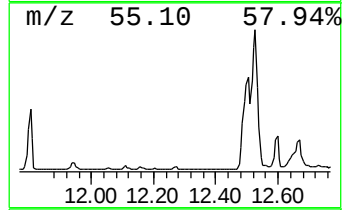
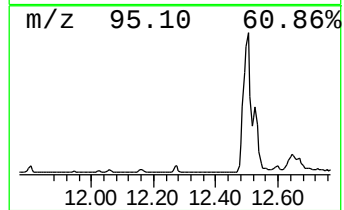
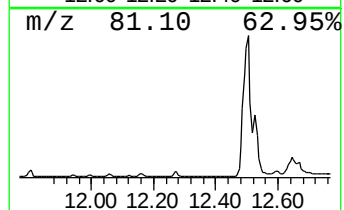
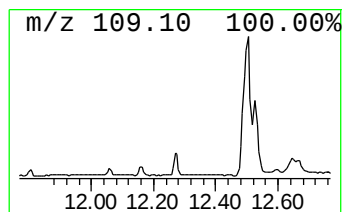
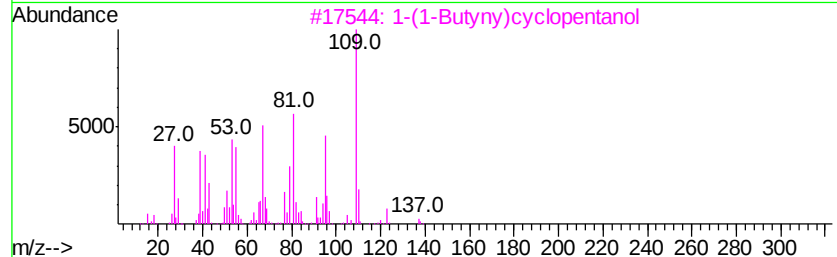
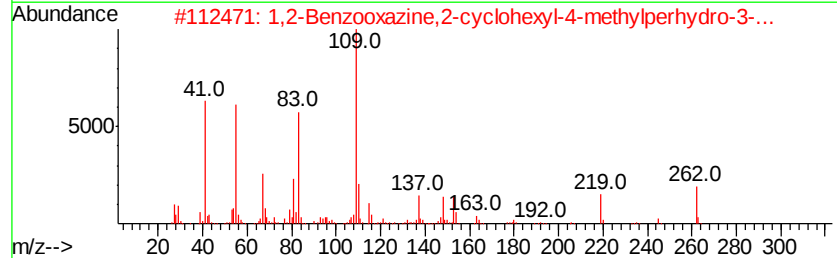
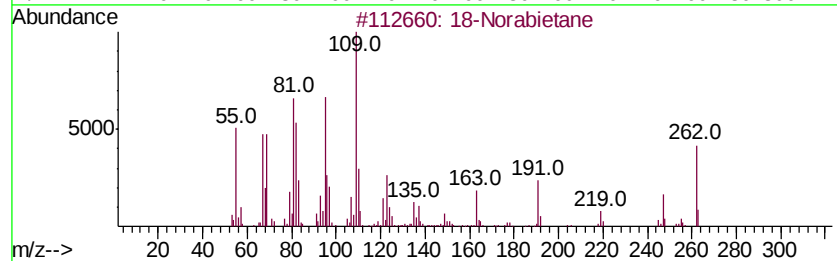
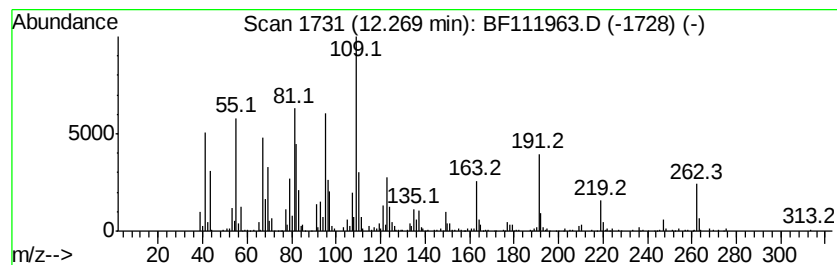
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 18-Norabietane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.27	10.07 ng	435604	Phenanthrene-d10	11.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	18-Norabietane	262	C19H34	1000293-16-6	94
2		1,2-Benzooxazine,2-cyclohexyl-4-...	262	C16H26N2O	037898-39-8	38
3		1-(1-Butynyl)cyclopentanol	138	C9H14O	1000342-86-5	35
4		8-Hexadecyne	222	C16H30	019781-86-3	22
5		Cyclopropane, 1,1-dimethyl-2-(2-...	124	C9H16	033422-32-1	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010919\
Data File : BF111963.D
Acq On : 9 Jan 2019 21:31
Operator : JU/SJ
Sample : K1044-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

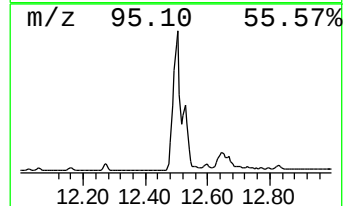
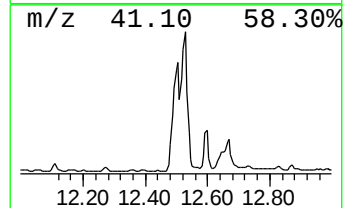
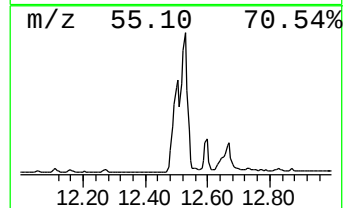
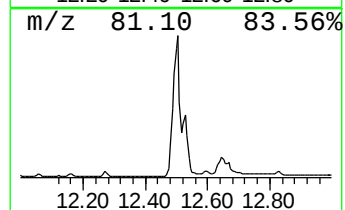
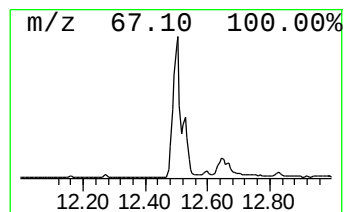
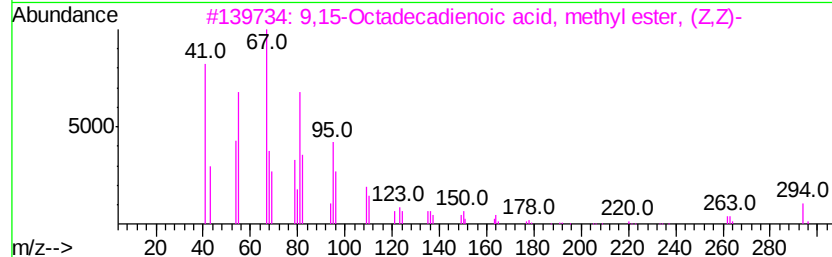
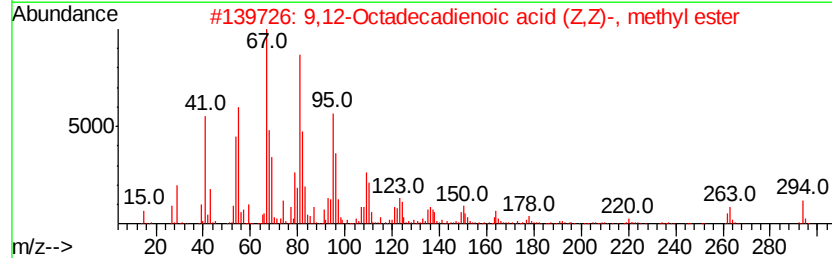
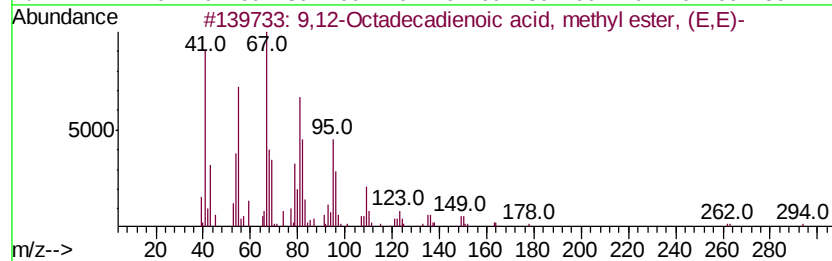
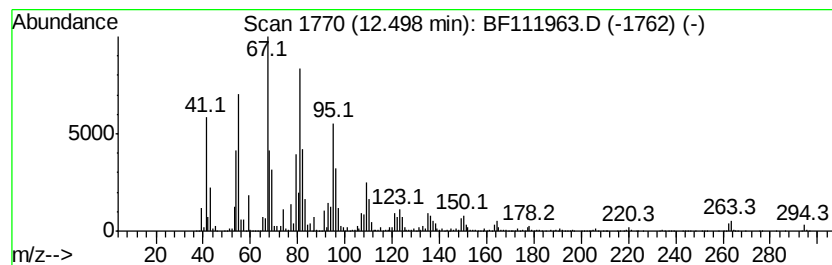
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 9,12-Octadecadienoic acid, ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.50	340.76 ng	14746500	Phenanthrene-d10	11.41	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
3	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
4	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
5	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010919\
Data File : BF111963.D
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ClientSampleId :
CPSB-11(20-25)

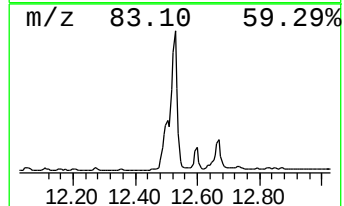
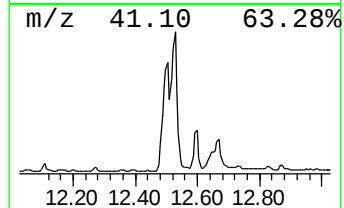
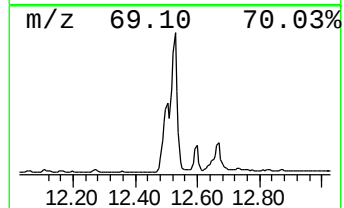
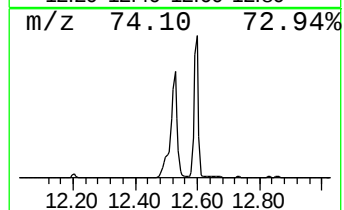
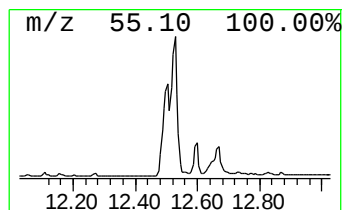
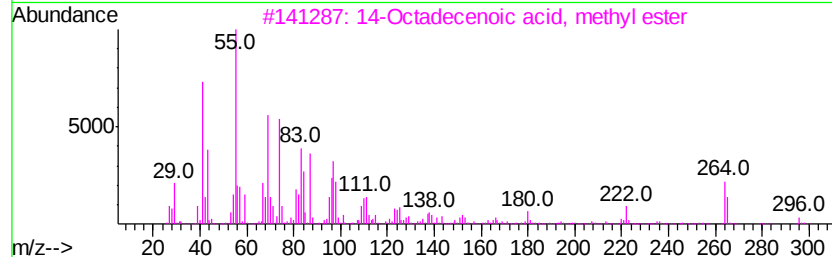
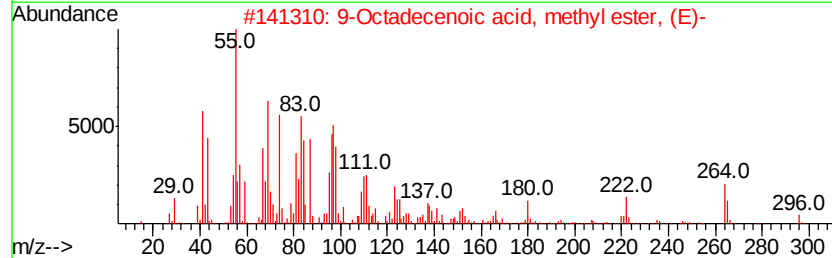
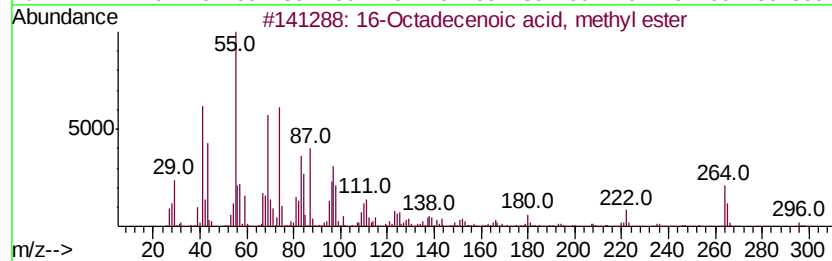
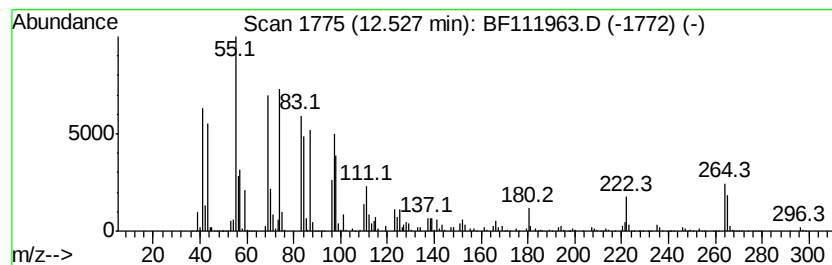
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 16-Octadecenoic acid, methyl ester Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.53	344.73 ng	14918300	Phenanthrene-d10	11.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	94
2		9-Octadecenoic acid, methyl ester...	296	C19H36O2	001937-62-8	90
3		14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	87
4		6-Octadecenoic acid, methyl ester...	296	C19H36O2	002777-58-4	86
5		8-Octadecenoic acid, methyl ester	296	C19H36O2	002345-29-1	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010919\
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Sample : K1044-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CPSB-11(20-25)

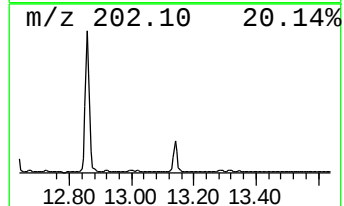
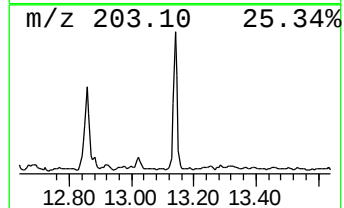
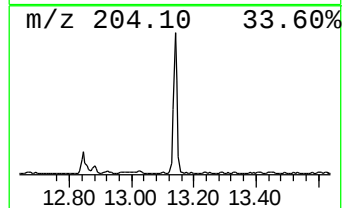
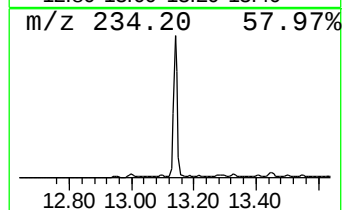
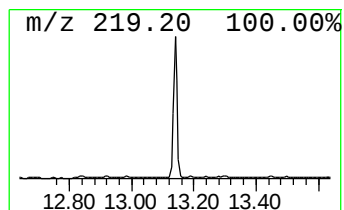
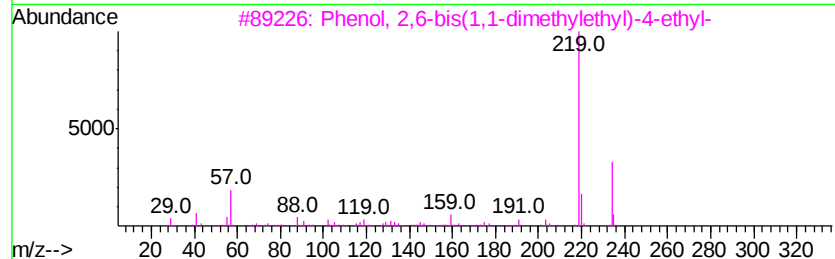
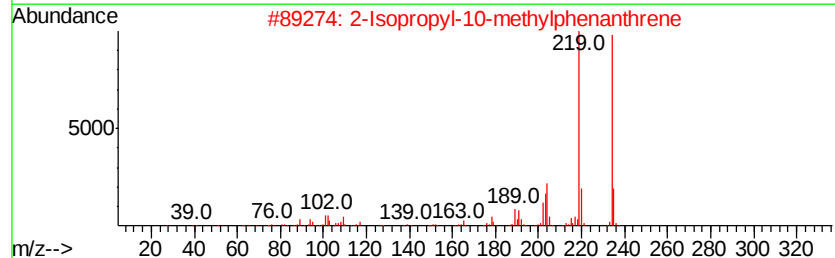
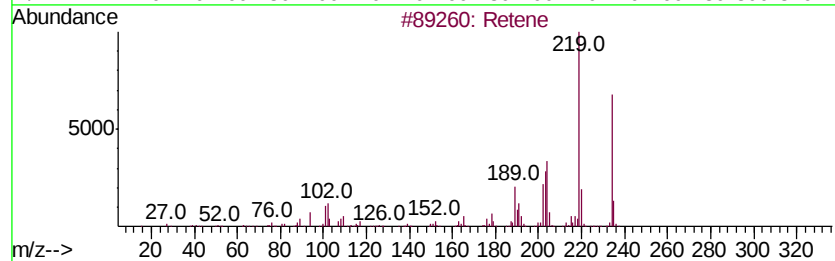
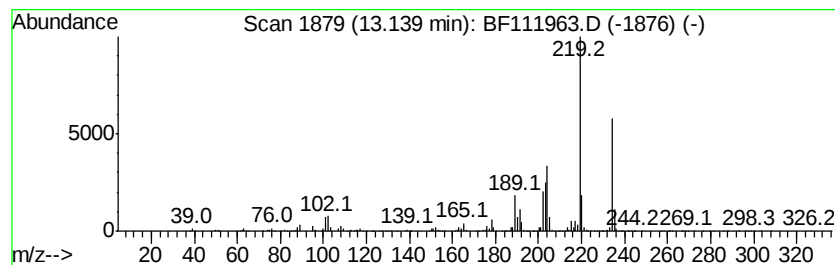
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Retene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.14	20.79 ng	879991	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Retene	234	C18H18	000483-65-8	98
2		2-Isopropyl-10-methylphenanthrene	234	C18H18	066552-97-4	95
3		Phenol, 2,6-bis(1,1-dimethylethyl)-	234	C16H26O	004130-42-1	52
4		1-Ethanone, 1-(4-amino-7-chloro-...	234	C12H11ClN2O	1000350-06-0	52
5		Ethanone, 1-(4,6-dihydroxy-2,3,5...	234	C13H14O4	021987-07-5	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010919\
Data File : BF111963.D
Acq On : 9 Jan 2019 21:31
Operator : JU/SJ
Sample : K1044-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CPSB-11(20-25)

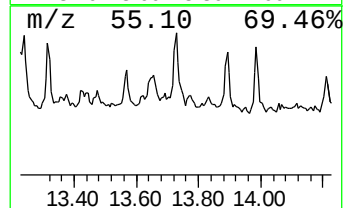
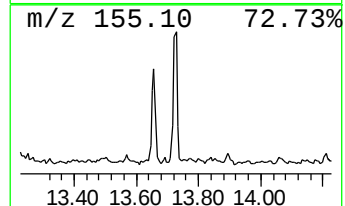
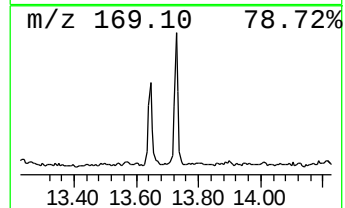
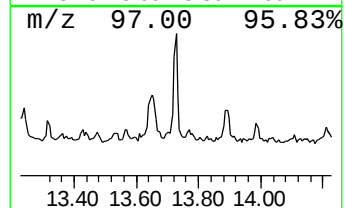
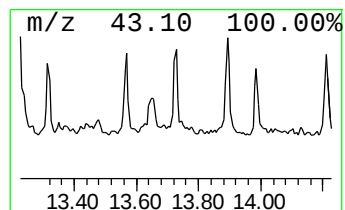
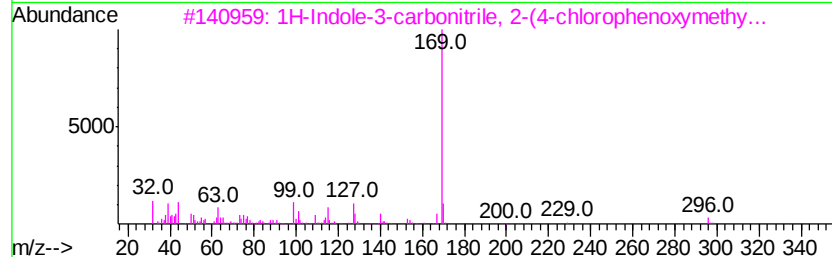
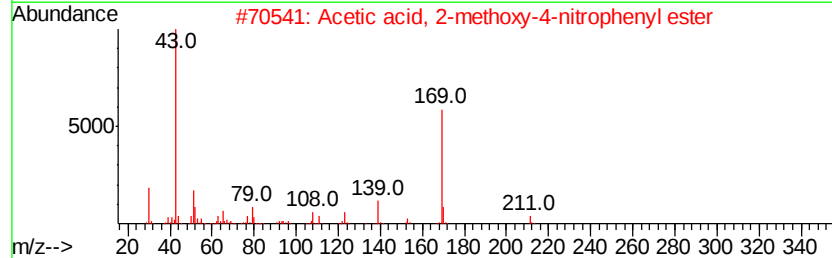
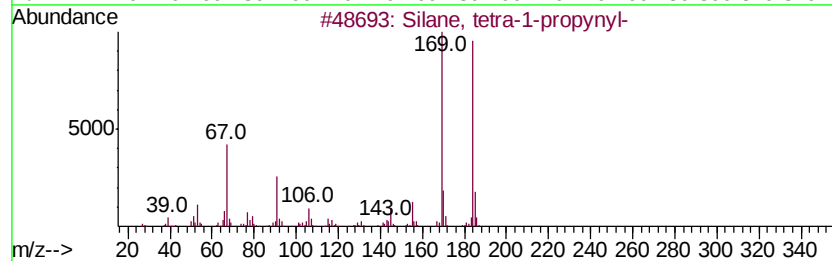
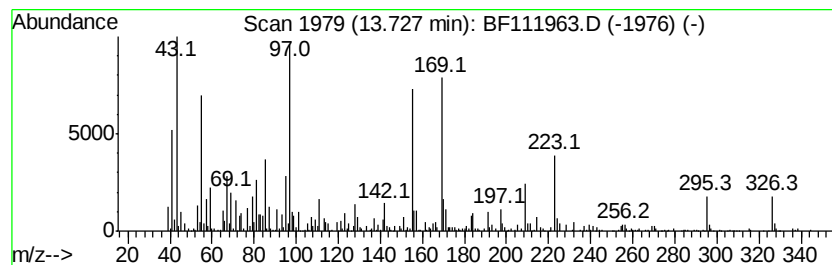
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown13.73 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	13.08 ng	553707	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Silane, tetra-1-propynyl-	184	C12H12Si	020143-21-9	22
2		Acetic acid, 2-methoxy-4-nitroph...	211	C9H9NO5	1000305-16-3	20
3		1H-Indole-3-carbonitrile, 2-(4-c...	296	C17H13ClN2O	1000304-36-8	18
4		1-methyl-1-cyclohexanecarboxylic...	224	C10H15F3O2	1000376-54-0	18
5		Ethanol, 1-(2-methyl-2H-tetrazol...	239	C9H13N5OS	1000316-81-4	15



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010919\
Data File : BF111963.D
Acq On : 9 Jan 2019 21:31
Operator : JU/SJ
Sample : K1044-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CPSB-11(20-25)

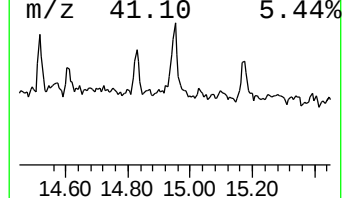
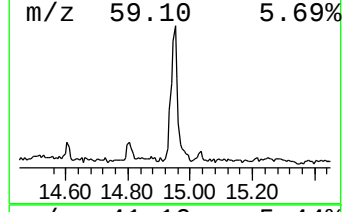
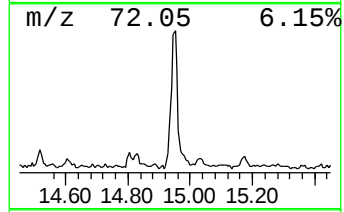
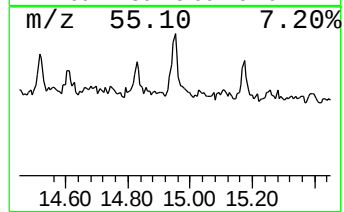
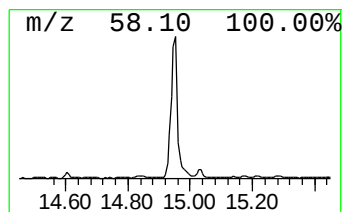
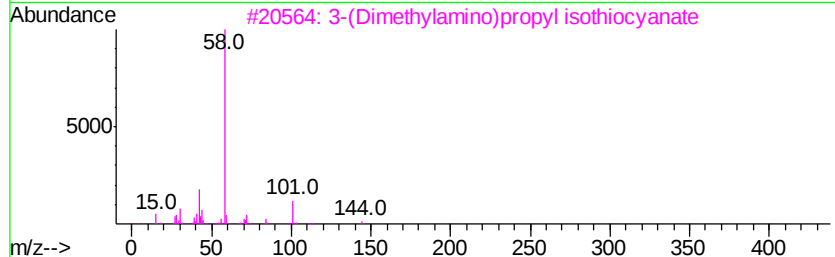
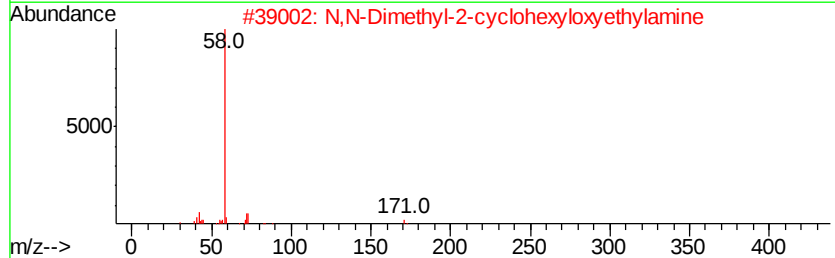
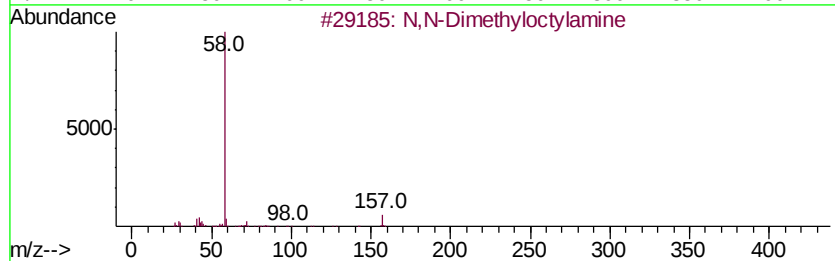
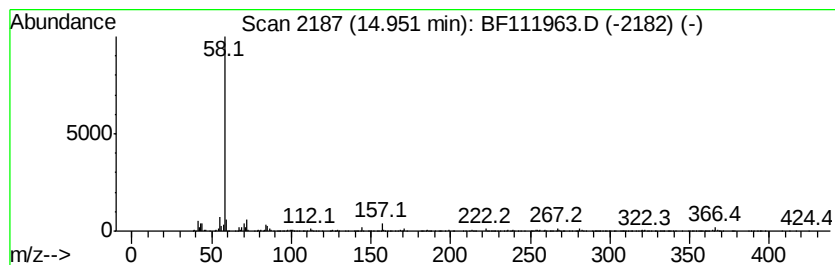
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 N,N-Dimethyloctylamine Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.95	20.82 ng	875864	Perylene-d12	15.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	N,N-Dimethyloctylamine	157	C10H23N	007378-99-6	80
2		N,N-Dimethyl-2-cyclohexyloxvethy...	171	C10H21NO	071126-60-8	64
3		3-(Dimethylamino)propyl isothiocy...	144	C6H12N2S	027421-70-1	64
4		2-Butanone, 4-(dimethylamino)-3-...	129	C7H15NO	022104-62-7	64
5		1-Hexanamine, 2-ethyl-N,N-dimethyl-	157	C10H23N	028056-87-3	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF010919\
 Data File : BF111963.D
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 Operator : JU/SJ
 Sample : K1044-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CPSB-11(20-25)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.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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Hexadecane	9.85	22.6	ng	1088170	3	9.92	964298	20.0
Heptacosane	10.57	10.5	ng	508332	3	9.92	964298	20.0
Tetracosane	11.27	15.8	ng	684471	4	11.41	865503	20.0
Bacchotricuneatin c	11.30	10.2	ng	439587	4	11.41	865503	20.0
Hexadecanoic acid...	11.80	143.0	ng	6187650	4	11.41	865503	20.0
n-Hexadecanoic acid	11.95	20.5	ng	887014	4	11.41	865503	20.0
18-Norabietane	12.27	10.1	ng	435604	4	11.41	865503	20.0
9,12-Octadecadien...	12.50	340.8	ng	14746500	4	11.41	865503	20.0
16-Octadecenoic a...	12.53	344.7	ng	14918300	4	11.41	865503	20.0
Retene	13.14	20.8	ng	879991	5	14.06	846751	20.0
unknown13.73	13.73	13.1	ng	553707	5	14.06	846751	20.0
N,N-Dimethyloctyl...	14.95	20.8	ng	875864	6	15.54	841214	20.0