

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
 Data File : BF102522.D
 Acq On : 27 Jan 2018 20:42
 Operator : SJ/JU
 Sample : J1009-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-06-012518-B

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.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	4.922	478	482	491	rVB	94707	143223	2.41%	0.218%
2	5.334	547	552	555	rBV	3608042	3605428	60.67%	5.478%
3	6.363	722	727	738	rBV	3100715	3847784	64.75%	5.846%
4	6.487	744	748	752	rBV	3720179	3719827	62.60%	5.652%
5	6.545	754	758	763	rVB2	80836	92679	1.56%	0.141%
6	6.716	783	787	793	rVB	1162915	1009341	16.99%	1.534%
7	6.869	808	813	822	rBV	3071572	2872840	48.35%	4.365%
8	7.104	849	853	856	rBV3	76540	90442	1.52%	0.137%
9	7.251	871	878	879	rBV2	68652	97211	1.64%	0.148%
10	7.287	879	884	886	rVV	2396940	2299450	38.70%	3.494%
11	7.304	886	887	891	rVB	396381	320269	5.39%	0.487%
12	7.357	891	896	899	rVB5	67055	96972	1.63%	0.147%
13	7.516	920	923	926	rVB2	138318	127043	2.14%	0.193%
14	7.610	932	939	941	rBV5	75199	116921	1.97%	0.178%
15	7.634	941	943	947	rVV4	115476	126175	2.12%	0.192%
16	7.740	958	961	967	rVV2	164501	202801	3.41%	0.308%
17	7.863	979	982	986	rBV3	66075	79656	1.34%	0.121%
18	7.945	991	996	998	rBV4	72185	106751	1.80%	0.162%
19	7.975	998	1001	1003	rVV	896175	899371	15.14%	1.367%
20	7.998	1003	1005	1008	rVV	1570555	1285998	21.64%	1.954%
21	8.051	1010	1014	1017	rVB	323425	365444	6.15%	0.555%
22	8.081	1017	1019	1023	rBV4	74762	73365	1.23%	0.111%
23	8.192	1036	1038	1041	rVB3	95759	81364	1.37%	0.124%
24	8.287	1048	1054	1060	rBV6	190791	295510	4.97%	0.449%
25	8.339	1060	1063	1065	rBV3	91739	72467	1.22%	0.110%
26	8.369	1065	1068	1072	rVB3	227127	218671	3.68%	0.332%
27	8.416	1072	1076	1078	rBV	687161	601902	10.13%	0.915%
28	8.498	1088	1090	1093	rVB2	115741	87844	1.48%	0.133%
29	8.587	1102	1105	1108	rBV	1155231	907002	15.26%	1.378%
30	8.645	1113	1115	1118	rVV3	98004	128203	2.16%	0.195%
31	8.681	1118	1121	1124	rVB3	160059	165426	2.78%	0.251%
32	8.710	1124	1126	1128	rBV	133054	92754	1.56%	0.141%
33	8.769	1133	1136	1139	rBV4	59044	86685	1.46%	0.132%
34	8.810	1140	1143	1146	rVB2	143305	161333	2.71%	0.245%

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
 Data File : BF102522.D
 Acq On : 27 Jan 2018 20:42
 Operator : SJ/JU
 Sample : J1009-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-06-012518-B

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	8.869	1151	1153	1155	rBV3	88081	88516	1.49%	0.134%
36	8.898	1155	1158	1163	rVB4	138582	199952	3.36%	0.304%
37	8.951	1163	1167	1170	rVB	177663	197629	3.33%	0.300%
38	8.986	1170	1173	1175	rBV2	120608	96451	1.62%	0.147%
39	9.016	1175	1178	1182	rVB	545497	469408	7.90%	0.713%
40	9.075	1182	1188	1191	rBV	3860279	3268905	55.01%	4.967%
41	9.151	1196	1201	1204	rBV2	1212419	1292235	21.75%	1.963%
42	9.269	1216	1221	1224	rVB4	68329	91536	1.54%	0.139%
43	9.339	1229	1233	1236	rBV	183051	221264	3.72%	0.336%
44	9.410	1242	1245	1247	rBV2	240415	265575	4.47%	0.404%
45	9.434	1247	1249	1251	rVV2	146899	162344	2.73%	0.247%
46	9.469	1251	1255	1257	rVV2	941501	906248	15.25%	1.377%
47	9.486	1257	1258	1262	rVB2	131436	111134	1.87%	0.169%
48	9.528	1262	1265	1269	rBV2	139309	101551	1.71%	0.154%
49	9.610	1276	1279	1281	rBV2	115891	111259	1.87%	0.169%
50	9.651	1284	1286	1288	rVV	134580	119145	2.01%	0.181%
51	9.681	1288	1291	1295	rVB	1005545	850711	14.32%	1.293%
52	9.728	1295	1299	1300	rBV2	97674	127994	2.15%	0.194%
53	9.751	1300	1303	1306	rVV	1496697	1248576	21.01%	1.897%
54	9.781	1306	1308	1312	rVB	240347	200961	3.38%	0.305%
55	9.851	1317	1320	1324	rVB	102084	120157	2.02%	0.183%
56	9.904	1324	1329	1332	rBV6	105005	150250	2.53%	0.228%
57	9.939	1332	1335	1336	rVV	94461	108740	1.83%	0.165%
58	9.951	1336	1337	1341	rVV3	147533	145379	2.45%	0.221%
59	9.992	1341	1344	1349	rVB3	187078	221622	3.73%	0.337%
60	10.063	1354	1356	1360	rBV	96153	114634	1.93%	0.174%
61	10.151	1368	1371	1373	rBV2	87413	118653	2.00%	0.180%
62	10.175	1373	1375	1378	rVB	763748	601566	10.12%	0.914%
63	10.298	1392	1396	1402	rVB2	176167	197122	3.32%	0.300%
64	10.392	1408	1412	1419	rVV	228950	306129	5.15%	0.465%
65	10.545	1434	1438	1441	rBV	1686628	1617691	27.22%	2.458%
66	10.651	1452	1456	1461	rBV2	613462	831119	13.99%	1.263%
67	10.839	1486	1488	1492	rBV5	58442	74099	1.25%	0.113%
68	11.098	1528	1532	1534	rBV	341607	302968	5.10%	0.460%
69	11.128	1534	1537	1543	rVB2	173199	222302	3.74%	0.338%
70	11.239	1552	1556	1558	rBV	1038920	1050404	17.68%	1.596%
71	11.263	1558	1560	1563	rVV	1102401	870503	14.65%	1.323%

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
 Data File : BF102522.D
 Acq On : 27 Jan 2018 20:42
 Operator : SJ/JU
 Sample : J1009-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-06-012518-B

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	11.310	1566	1568	1571	rVV	168148	164184	2.76%	0.249%
73	11.351	1571	1575	1580	rVB3	45477	73485	1.24%	0.112%
74	11.475	1591	1596	1601	rVB2	143851	152385	2.56%	0.232%
75	11.522	1601	1604	1607	rBV	226778	187204	3.15%	0.284%
76	11.628	1618	1622	1625	rBV	2745095	2214486	37.27%	3.365%
77	11.739	1636	1641	1643	rBV	112622	125052	2.10%	0.190%
78	11.769	1643	1646	1649	rVB	151985	135512	2.28%	0.206%
79	11.851	1656	1660	1662	rBV2	191807	185032	3.11%	0.281%
80	11.927	1670	1673	1678	rBV	158257	189575	3.19%	0.288%
81	12.033	1688	1691	1693	rBV	112424	89527	1.51%	0.136%
82	12.063	1693	1696	1702	rVB	109993	123762	2.08%	0.188%
83	12.286	1731	1734	1735	rBV	82514	74567	1.25%	0.113%
84	12.322	1735	1740	1741	rVV	5240903	5942305	100.00%	9.029%
85	12.339	1741	1743	1748	rVB2	4653790	4319716	72.69%	6.563%
86	12.416	1753	1756	1759	rVV	1010361	920339	15.49%	1.398%
87	12.451	1759	1762	1767	rVB	1395363	1203599	20.25%	1.829%
88	12.680	1795	1801	1807	rVB	1070886	1271535	21.40%	1.932%
89	12.822	1818	1825	1834	rBV	2285452	2350558	39.56%	3.571%
90	12.904	1834	1839	1842	rBV3	70278	132028	2.22%	0.201%
91	13.010	1850	1857	1860	rBV	233850	311521	5.24%	0.473%
92	13.045	1861	1863	1865	rBV	91254	84410	1.42%	0.128%
93	13.074	1865	1868	1871	rVB	174676	166992	2.81%	0.254%
94	13.627	1960	1962	1964	rBV	109044	84375	1.42%	0.128%
95	13.716	1974	1977	1982	rVB3	209402	310609	5.23%	0.472%
96	13.880	2000	2005	2007	rBV2	1030913	1274866	21.45%	1.937%
97	13.904	2007	2009	2013	rVB	572144	449508	7.56%	0.683%
98	14.904	2176	2179	2186	rVB	368740	563798	9.49%	0.857%
99	15.251	2235	2238	2244	rVB	295953	368723	6.21%	0.560%
100	15.316	2245	2249	2252	rBV	572539	681774	11.47%	1.036%

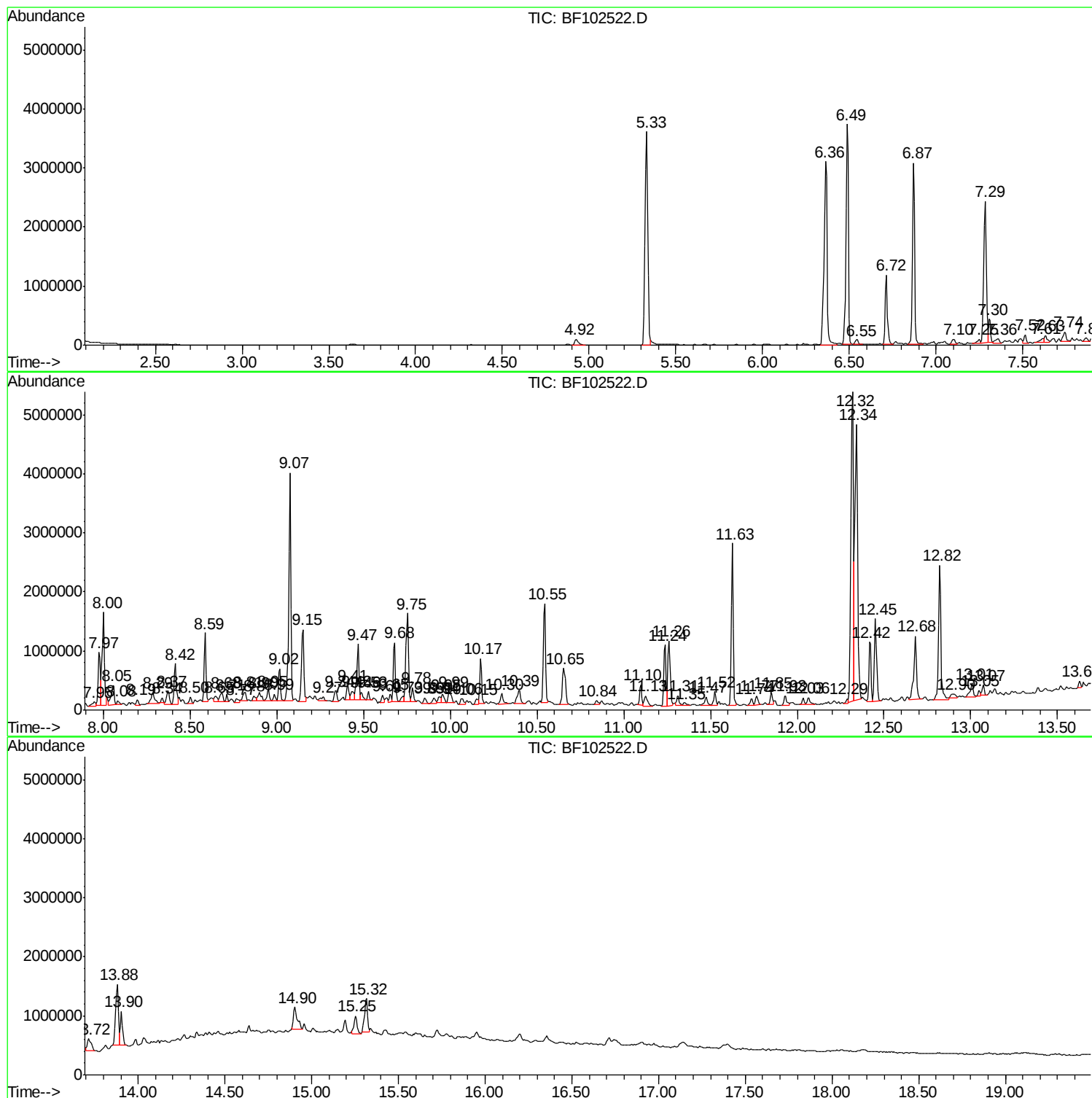
Sum of corrected areas: 65814311

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

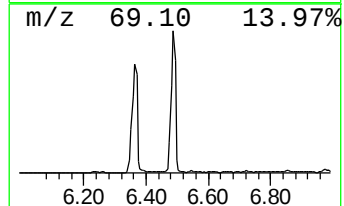
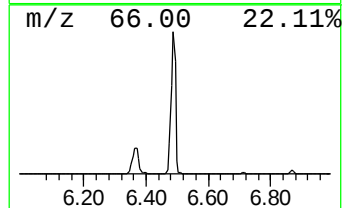
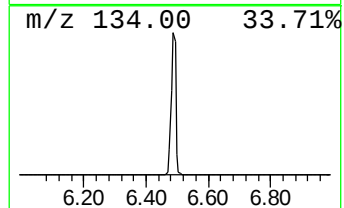
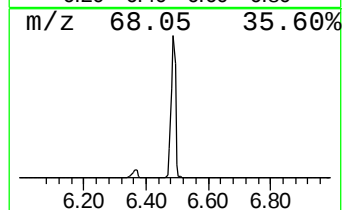
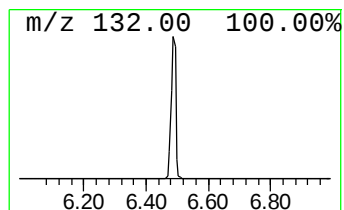
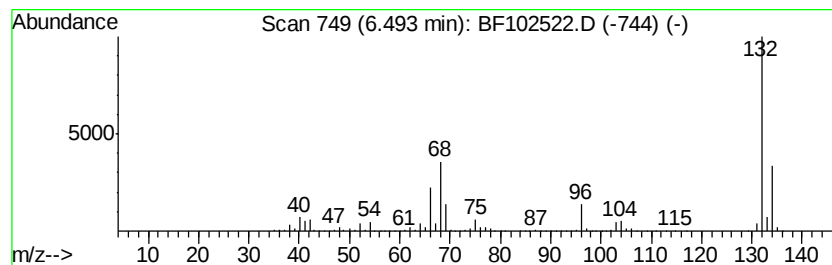
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.49 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	73.71 ng	3719830	1,4-Dichlorobenzene-d4	6.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

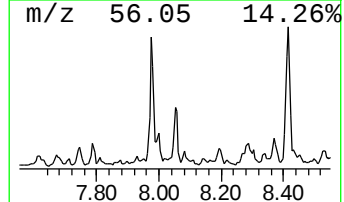
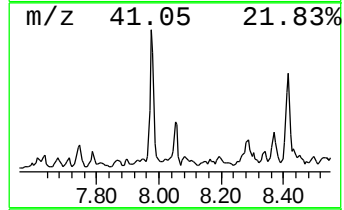
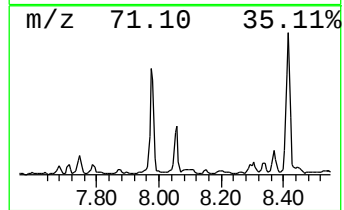
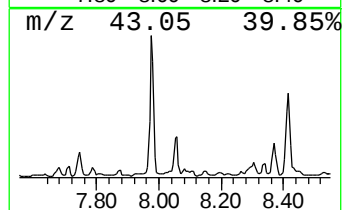
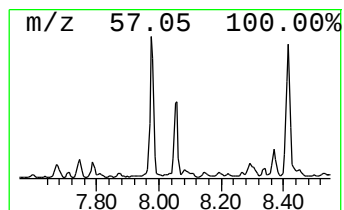
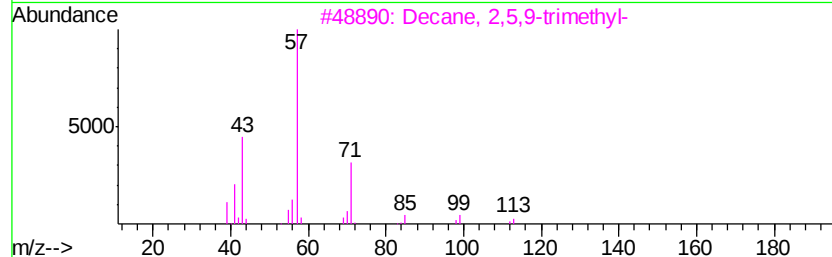
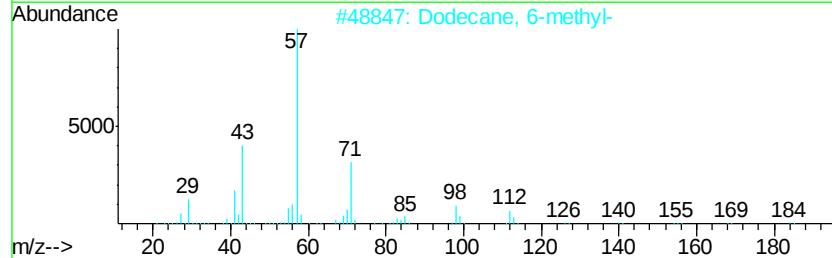
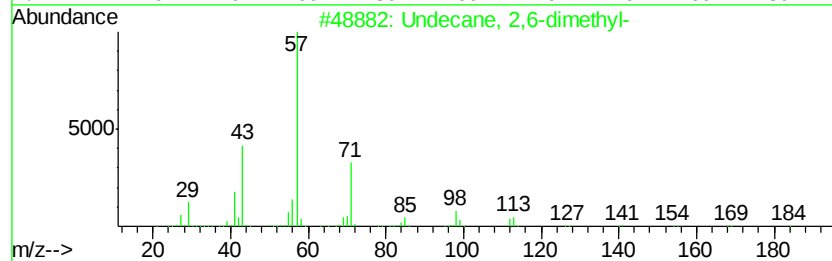
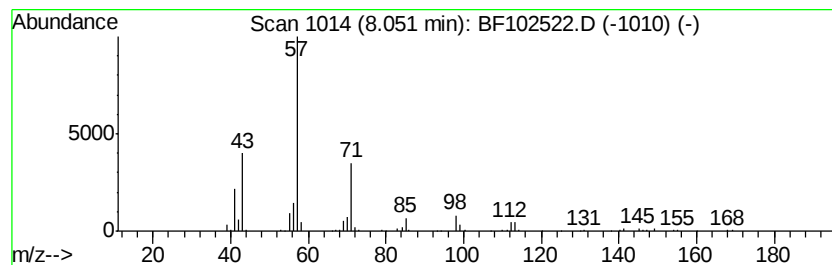
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Undecane, 2,6-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.05	5.68 ng	365444	Napthalene-d8	8.00

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	86	
2	Dodecane, 6-methyl-	184	C13H28	006044-71-9	72	
3	Decane, 2,5,9-trimethyl-	184	C13H28	062108-22-9	64	
4	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	59	
5	Octane, 2,5-dimethyl-	142	C10H22	015869-89-3	59	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleID :
AU-06-012518-B

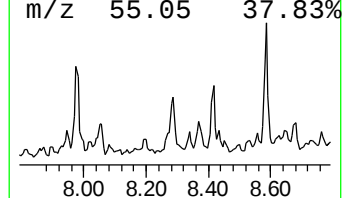
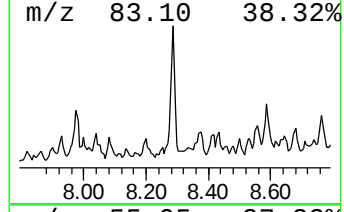
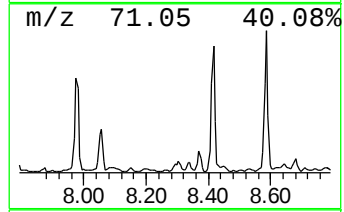
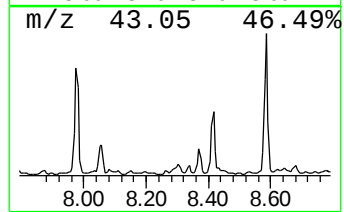
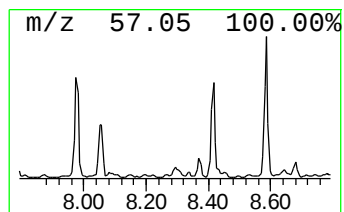
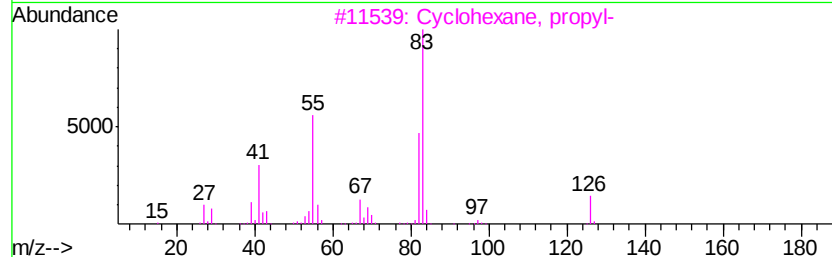
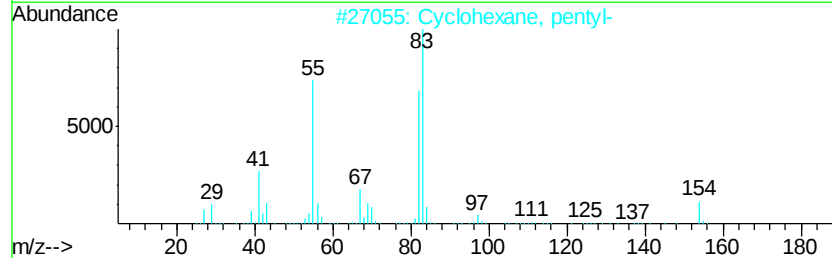
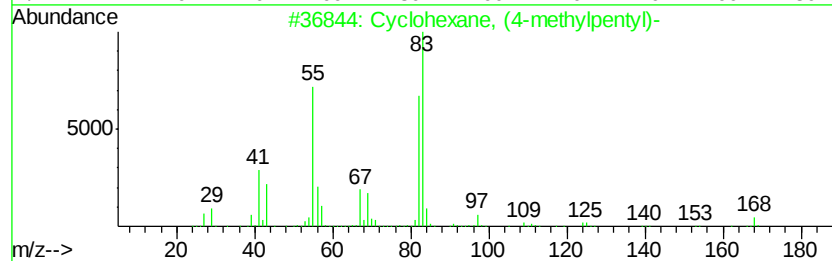
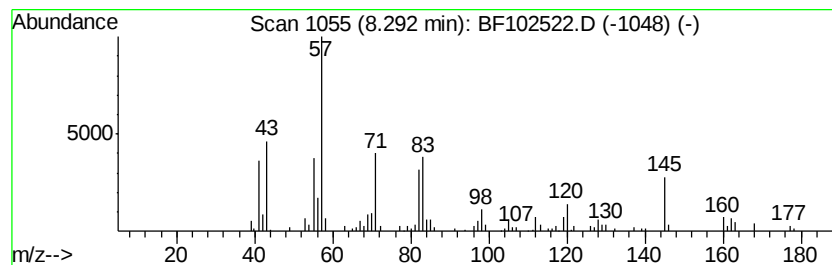
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Cyclohexane, (4-methylpentyl)- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.29	4.60 ng	295510	Napthalene-d8	8.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (4-methylpentyl)-	168	C12H24	061142-20-9	81
2			Cyclohexane, pentyl-	154	C11H22	004292-92-6	59
3			Cyclohexane, propyl-	126	C9H18	001678-92-8	59
4			Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	59
5			Cyclohexane, octyl-	196	C14H28	001795-15-9	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

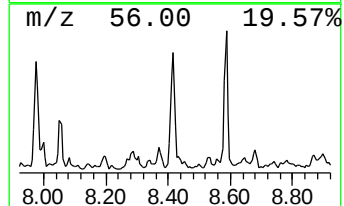
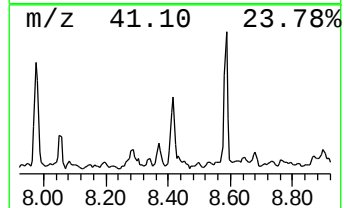
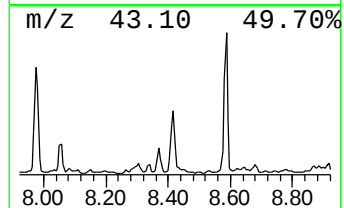
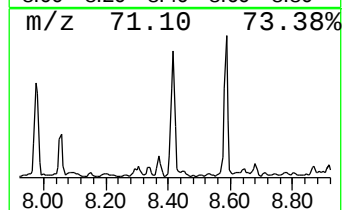
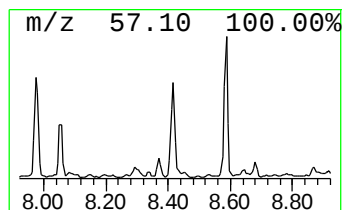
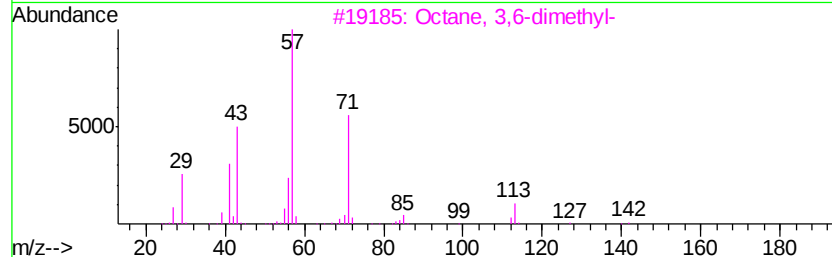
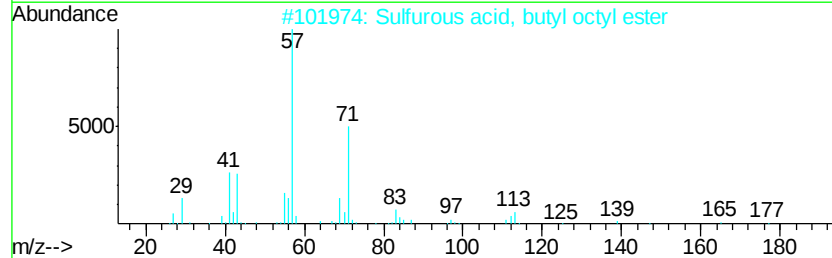
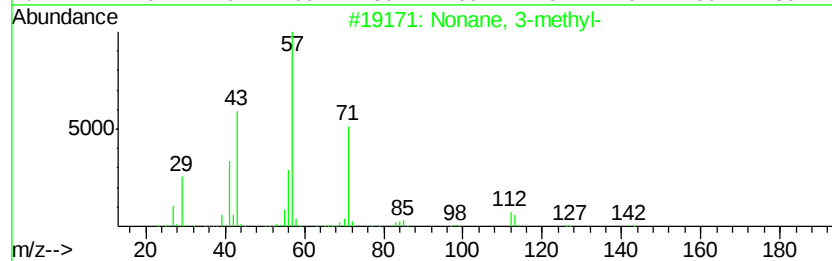
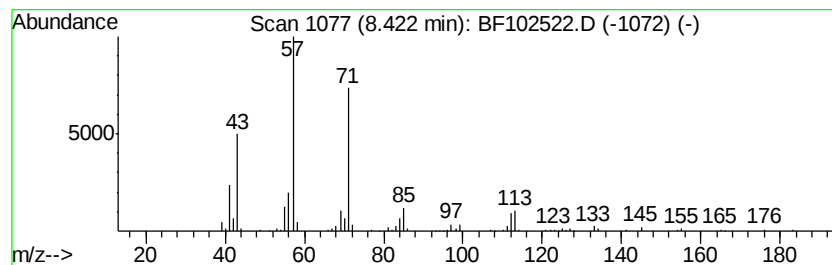
Instrument :
BNA_F
ClientSampled :
AU-06-012518-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Nonane, 3-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
8.42	9.36 ng	601902	Naphthalene-d8	8.00		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 3-methyl-	142	C10H22	005911-04-6	72	
2	Sulfurous acid, butyl octyl ester	250	C12H26O3S	1000309-17-5	64	
3	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	64	
4	Sulfurous acid, 2-ethylhexyl tri...	376	C21H44O3S	1000309-19-6	59	
5	Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	59	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

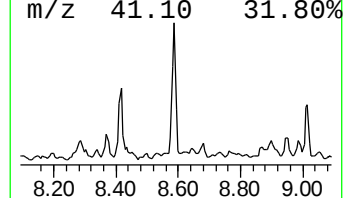
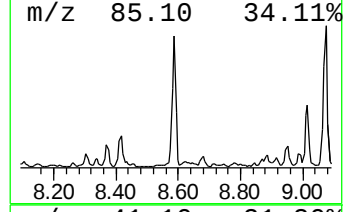
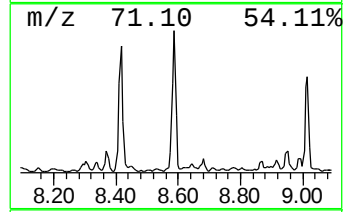
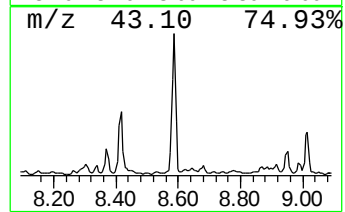
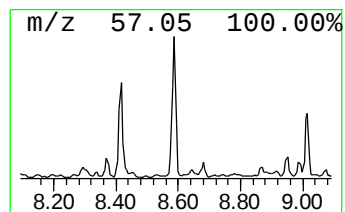
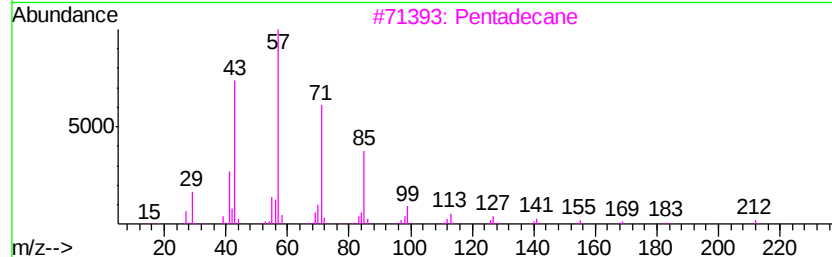
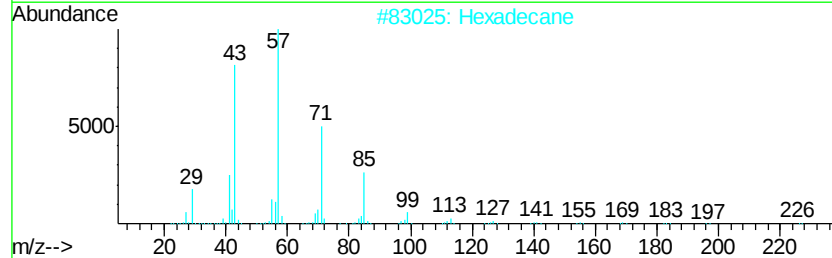
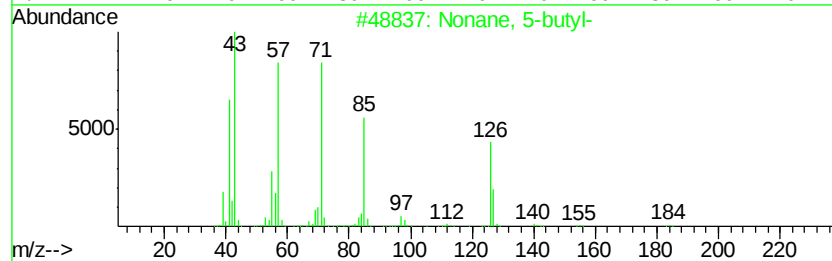
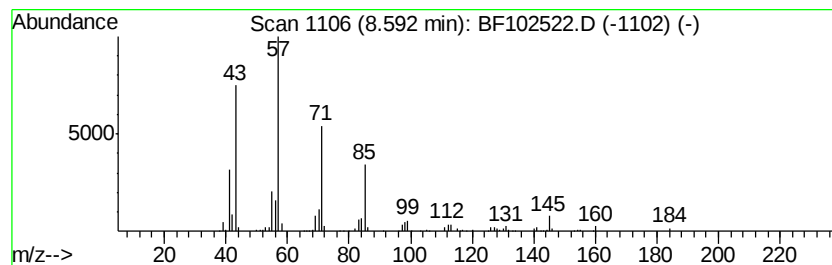
Instrument :
BNA_F
ClientSampleID :
AU-06-012518-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Nonane, 5-butyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.59	14.11 ng	907002	Naphthalene-d8	8.00
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Nonane, 5-butyl-	184 C13H28	017312-63-9	90
2	Hexadecane	226 C16H34	000544-76-3	90
3	Pentadecane	212 C15H32	000629-62-9	90
4	Tetradecane	198 C14H30	000629-59-4	90
5	Tridecane	184 C13H28	000629-50-5	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

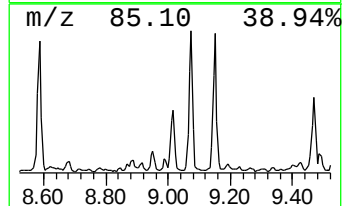
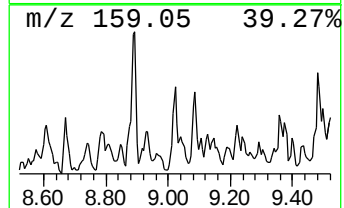
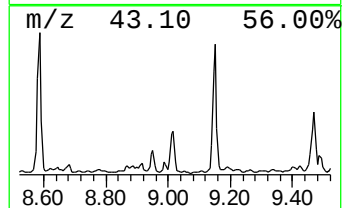
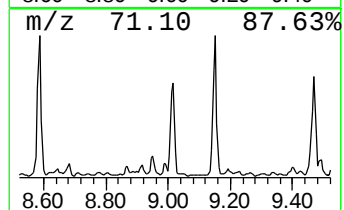
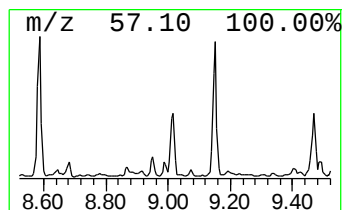
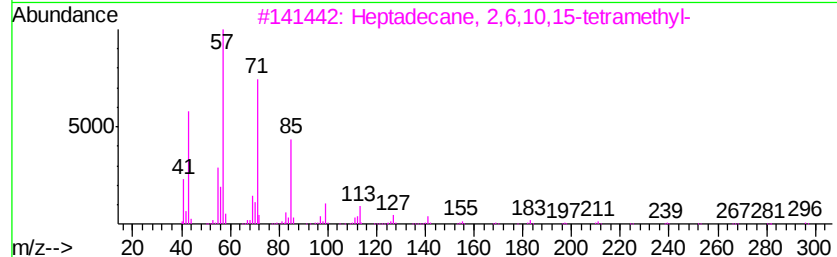
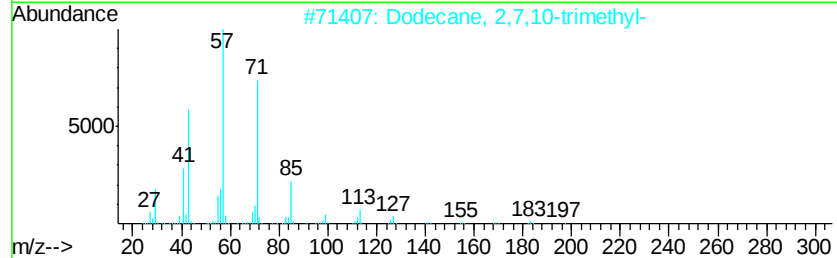
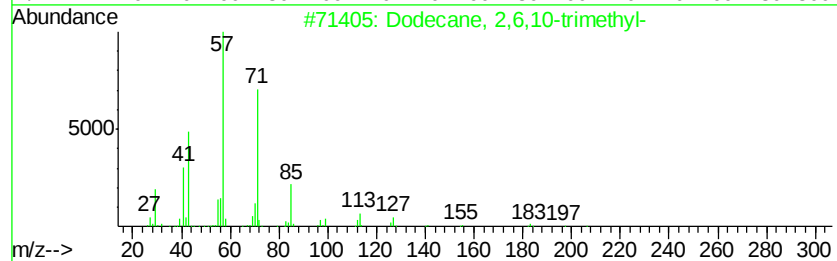
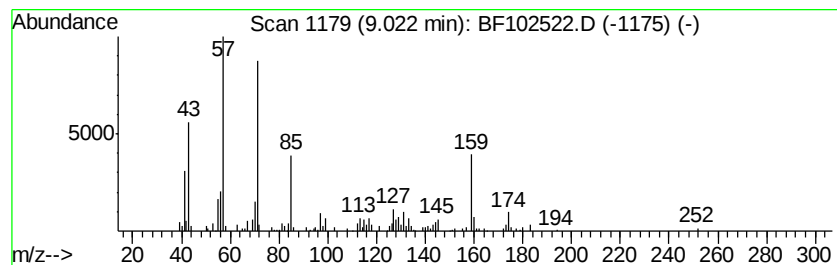
Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Dodecane, 2,6,10-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.02	7.52 ng	469408	Acenaphthene-d10	9.75	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	90
2	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	90
3	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	78
4	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
5	Bacchotricuneatin c	342	C20H22O5	066563-30-2	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

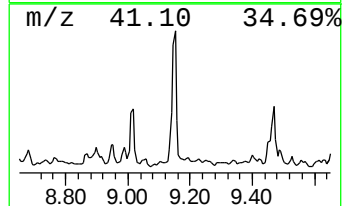
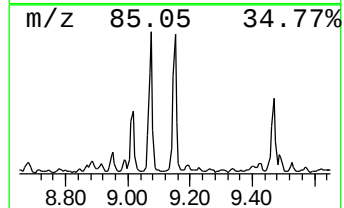
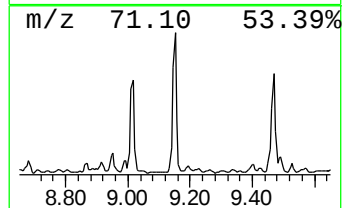
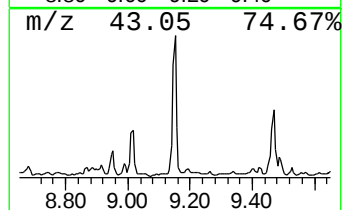
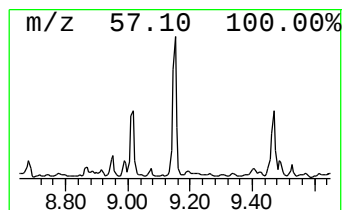
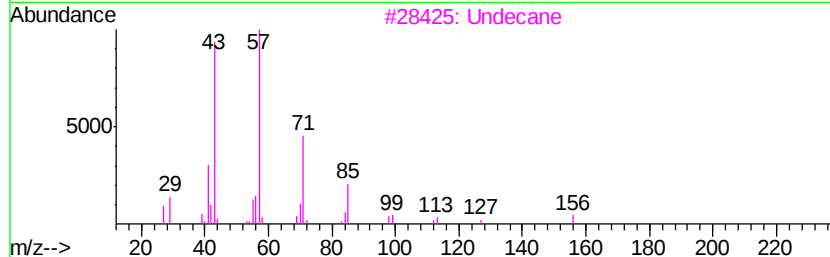
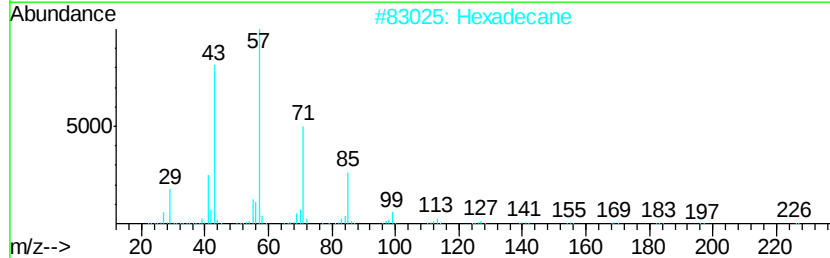
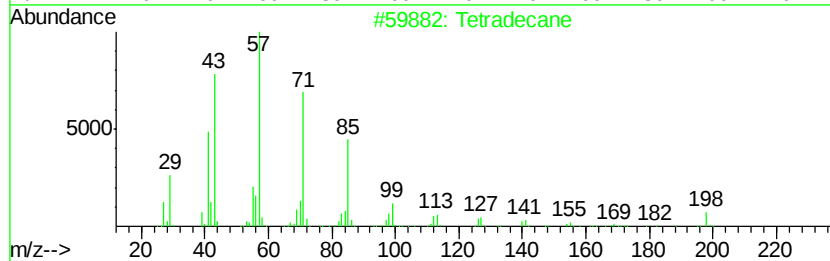
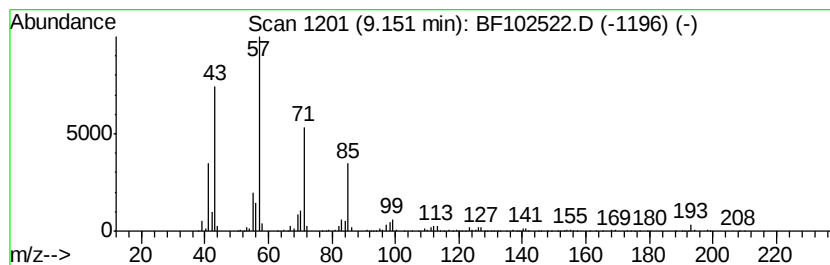
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tetradecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.15	20.70 ng	1292240	Acenaphthene-d10	9.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	94
2		Hexadecane	226	C16H34	000544-76-3	90
3		Undecane	156	C11H24	001120-21-4	86
4		Tridecane, 6-methyl-	198	C14H30	013287-21-3	74
5		Pentadecane	212	C15H32	000629-62-9	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

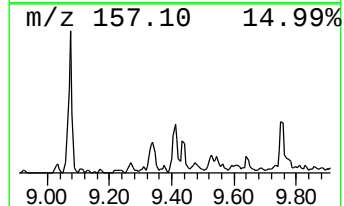
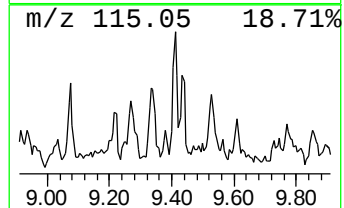
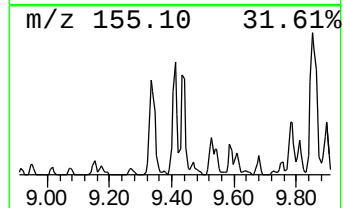
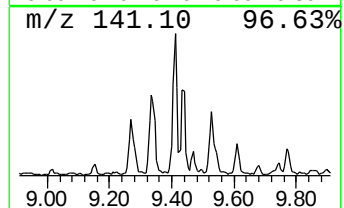
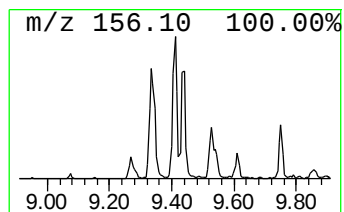
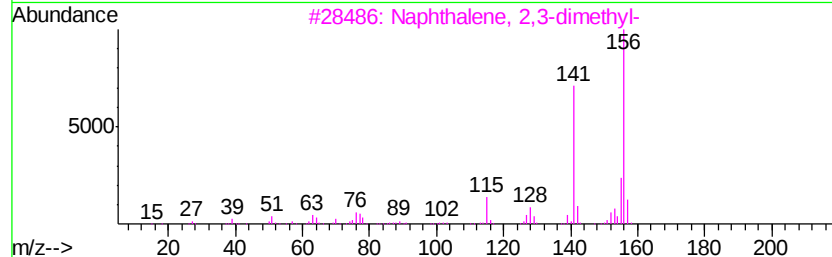
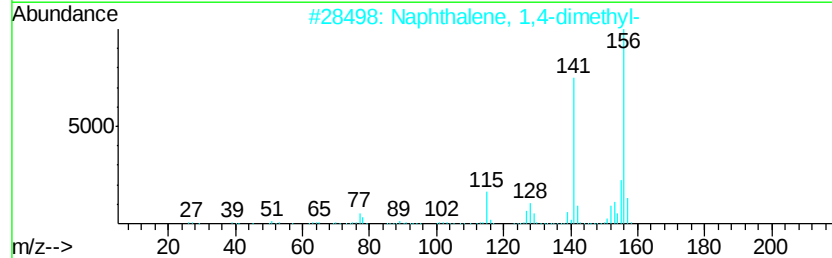
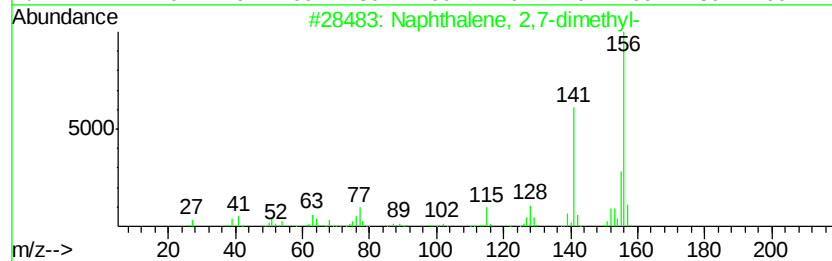
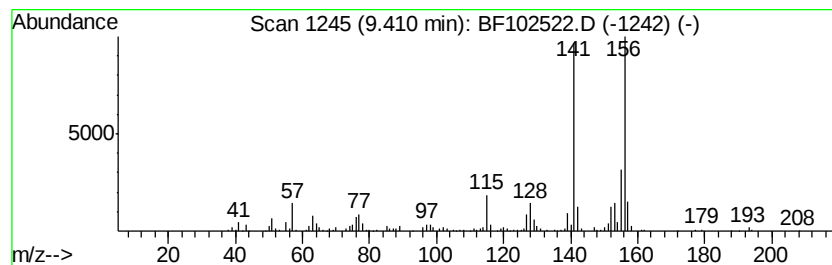
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Naphthalene, 2,7-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.41	4.25 ng	265575	Acenaphthene-d10	9.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
2		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
3		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
5		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

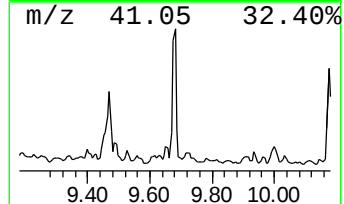
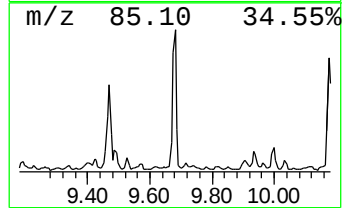
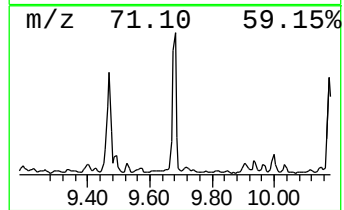
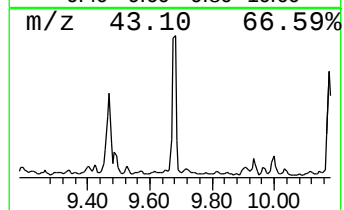
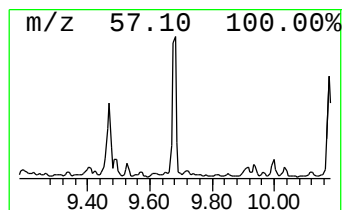
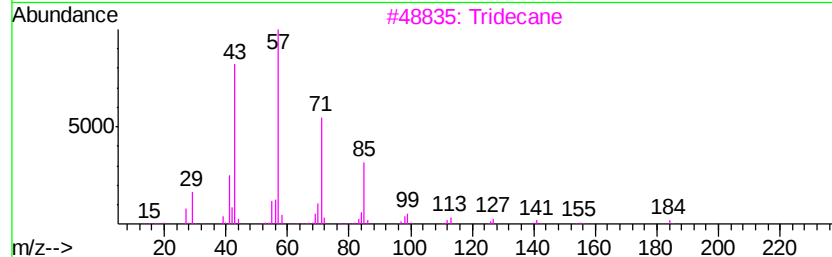
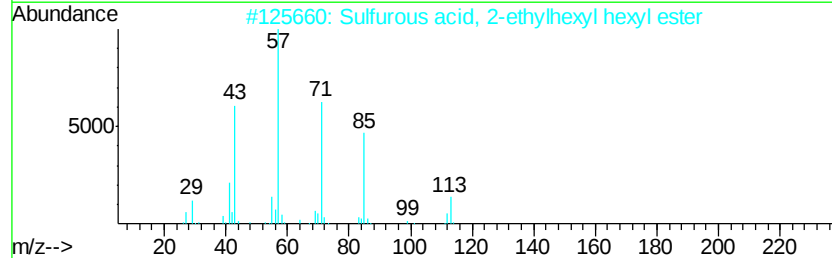
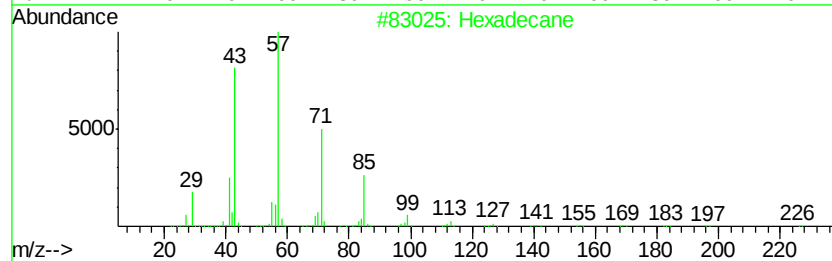
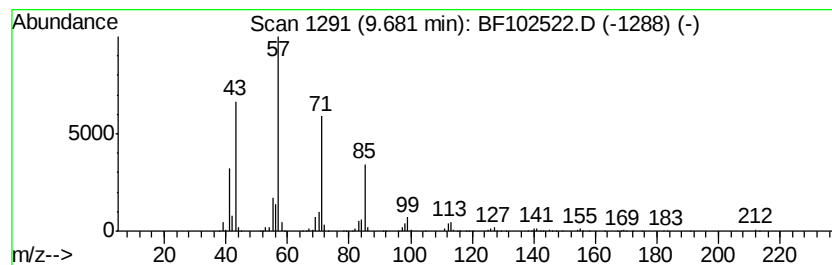
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.68	13.63 ng	850711	Acenaphthene-d10	9.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	90
2		Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	86
3		Tridecane	184	C13H28	000629-50-5	83
4		Tetradecane	198	C14H30	000629-59-4	64
5		Undecane, 6-ethyl-	184	C13H28	017312-60-6	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

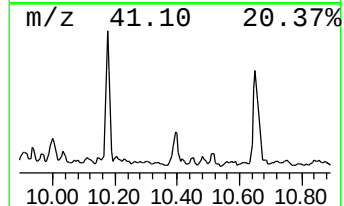
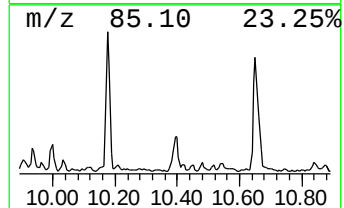
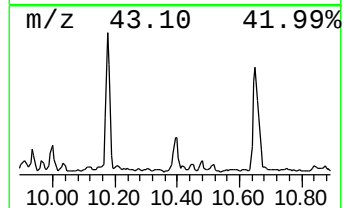
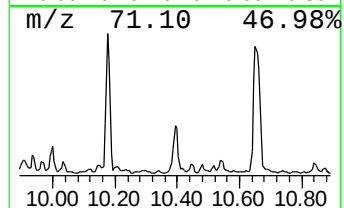
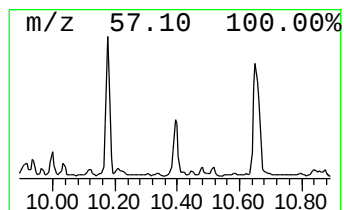
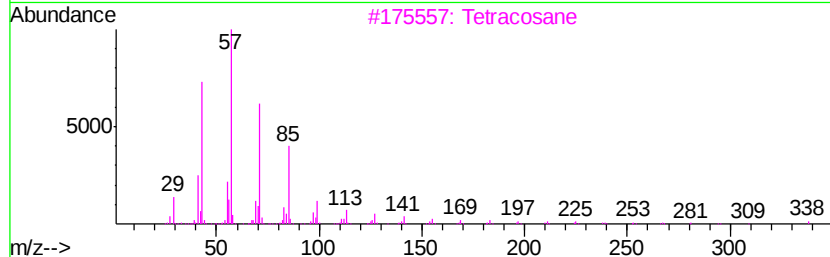
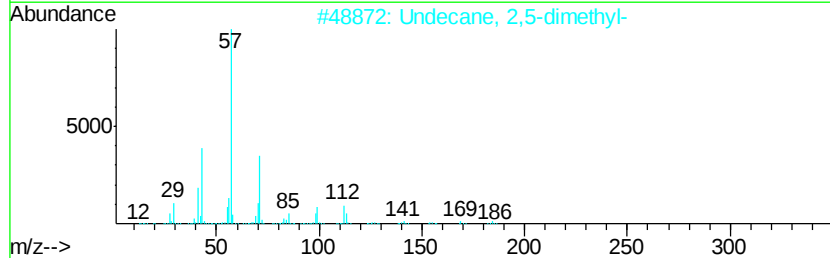
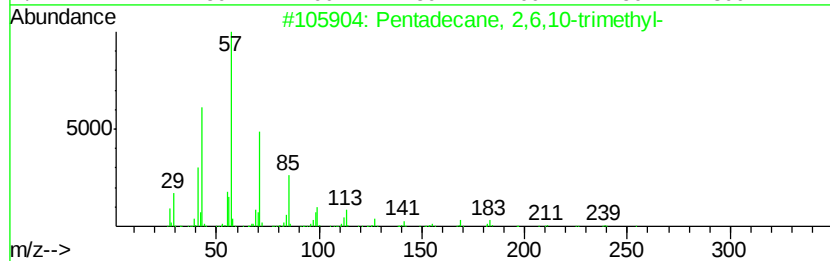
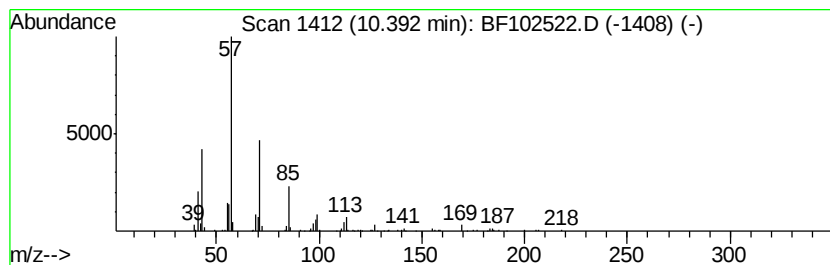
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pentadecane, 2,6,10-trimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.39	4.90 ng	306129	Acenaphthene-d10	9.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	90
2		Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	76
3		Tetracosane	338	C24H50	000646-31-1	72
4		Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	68
5		Pentadecane, 7-methyl-	226	C16H34	006165-40-8	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

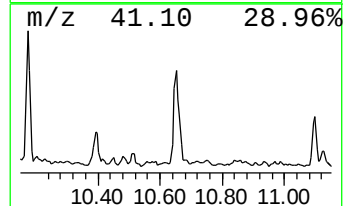
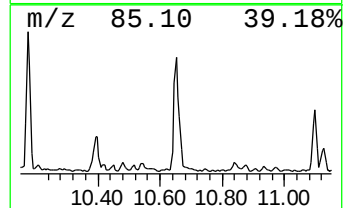
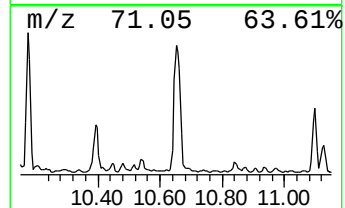
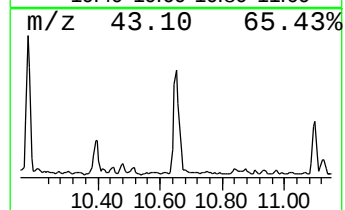
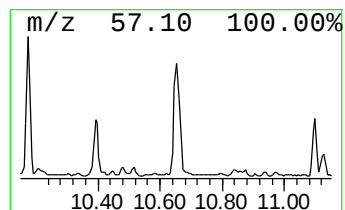
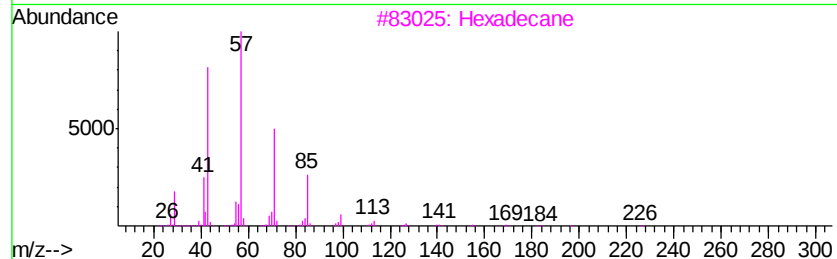
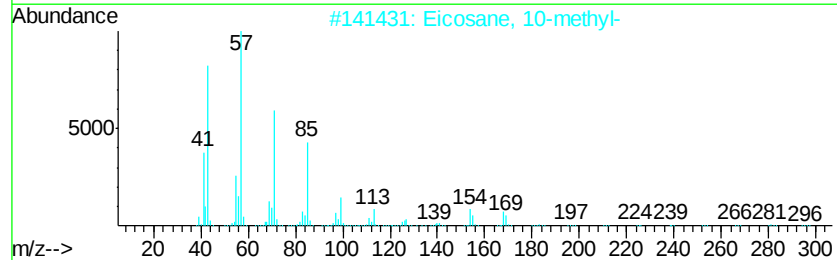
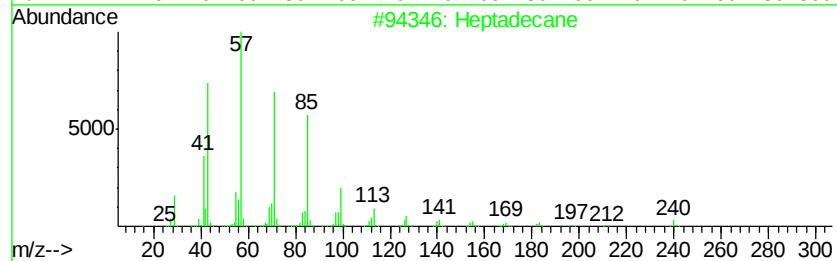
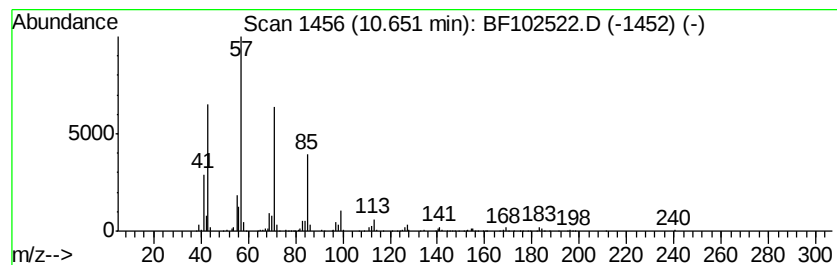
Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.65	15.82 ng	831119	Phenanthrene-d10		11.24
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	97
2	Eicosane, 10-methyl-	296	C21H44	054833-23-7	90
3	Hexadecane	226	C16H34	000544-76-3	90
4	Tetracosane	338	C24H50	000646-31-1	90
5	Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

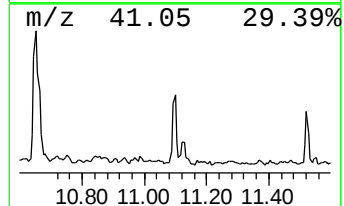
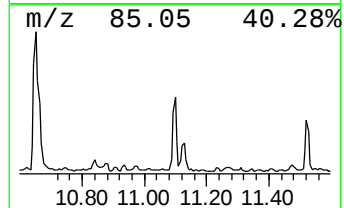
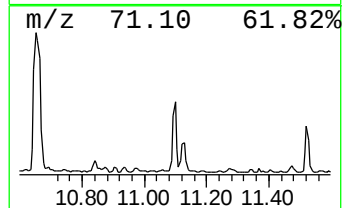
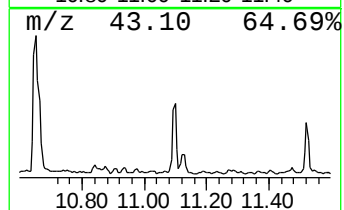
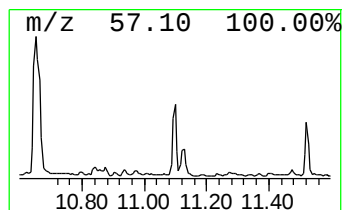
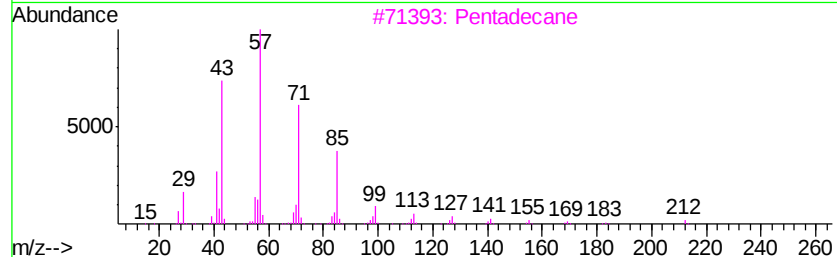
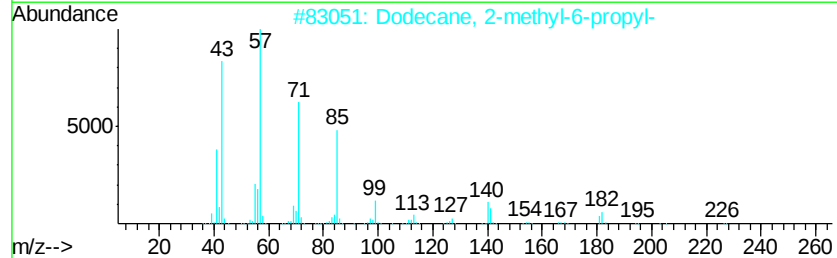
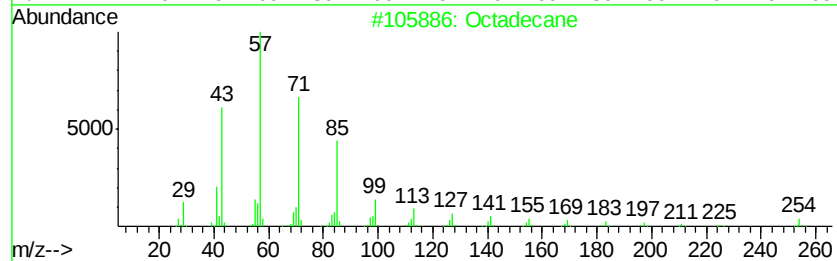
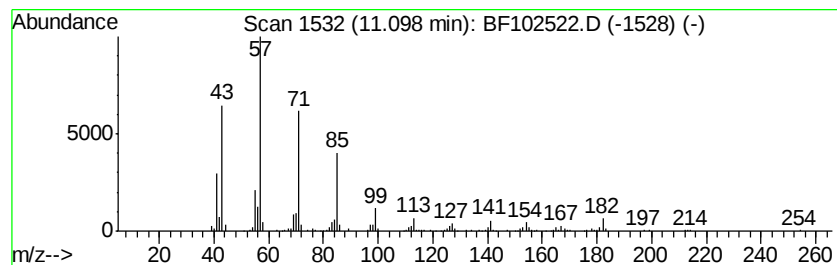
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Octadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.10	5.77 ng	302968	Phenanthrene-d10	11.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	94
2		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90
3		Pentadecane	212	C15H32	000629-62-9	87
4		Hexadecane	226	C16H34	000544-76-3	87
5		Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

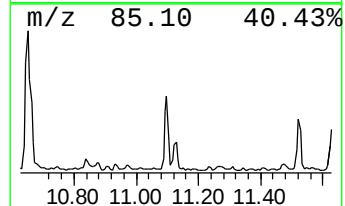
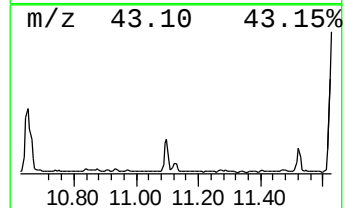
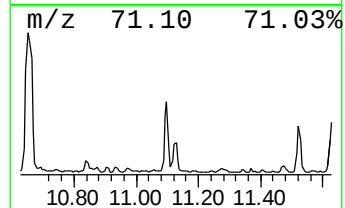
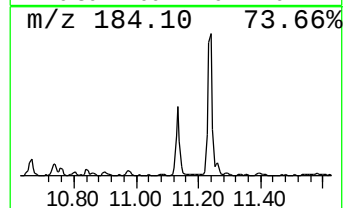
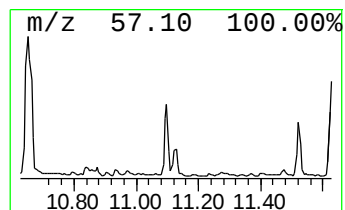
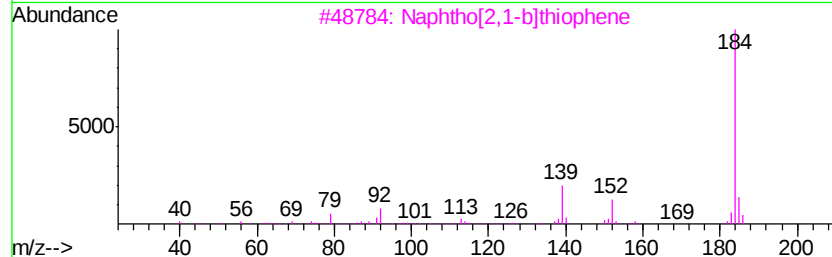
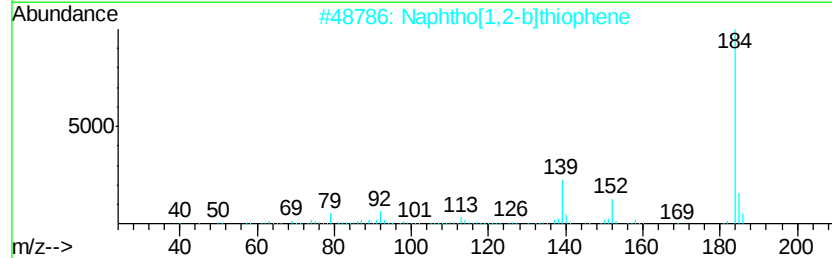
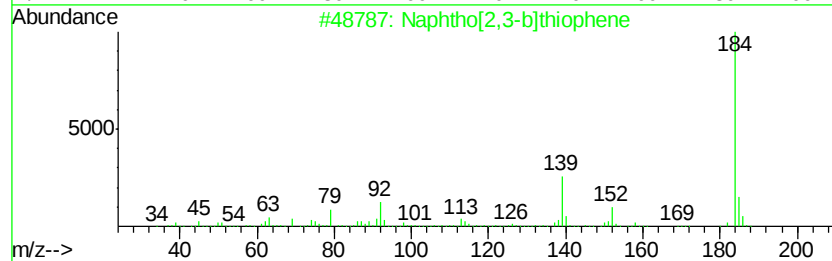
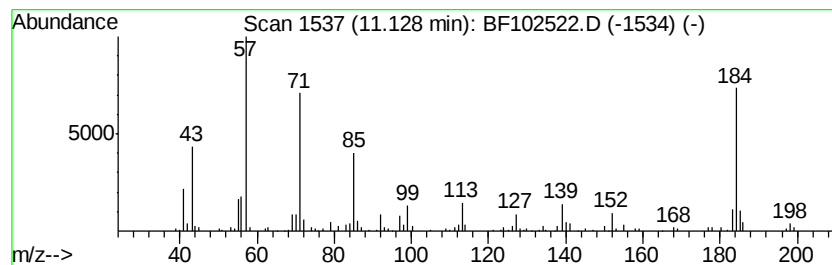
Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Naphtho[2,3-b]thiophene Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.13	4.23 ng	222302	Phenanthrene-d10		11.24	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	55
2		Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	55
3		Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	55
4		Dibenzothiophene	184	C12H8S	000132-65-0	55
5		Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	49



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

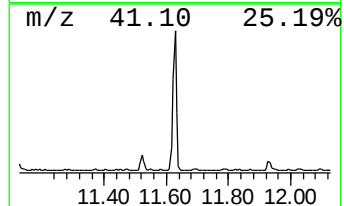
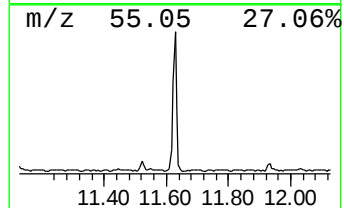
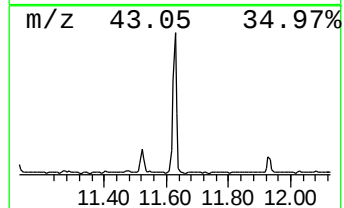
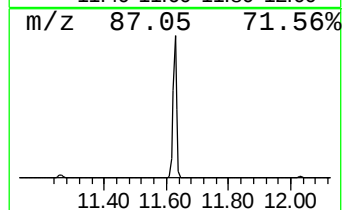
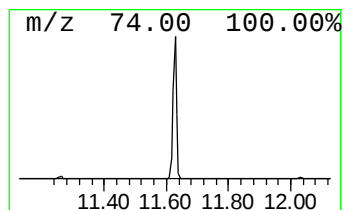
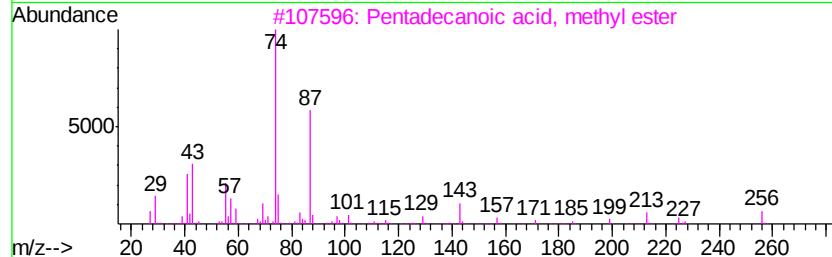
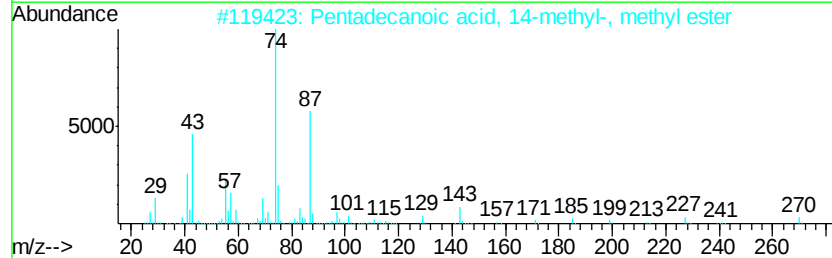
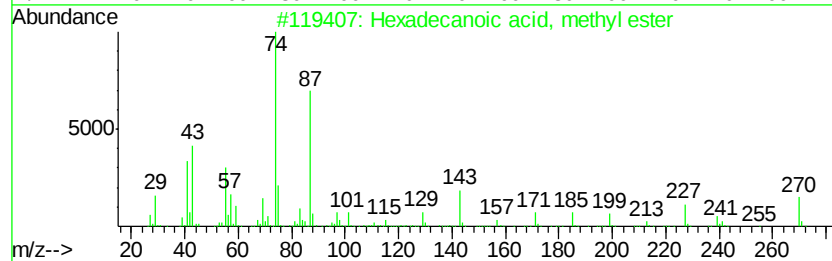
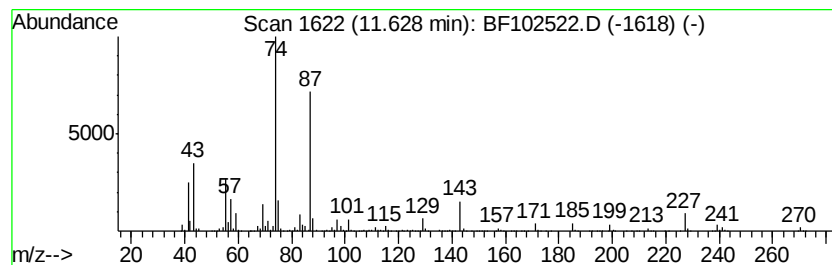
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Hexadecanoic acid, methyl e... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.63	42.16 ng	2214490	Phenanthrene-d10	11.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	93
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	91
3		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	90
4		Methyl tetradecanoate	242	C15H30O2	000124-10-7	80
5		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

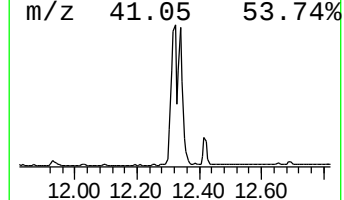
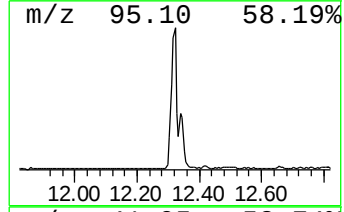
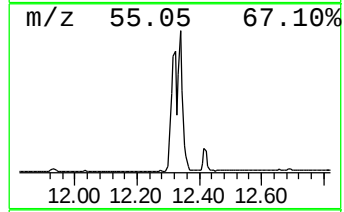
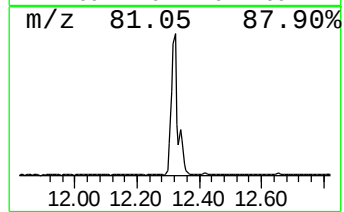
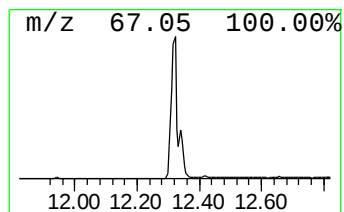
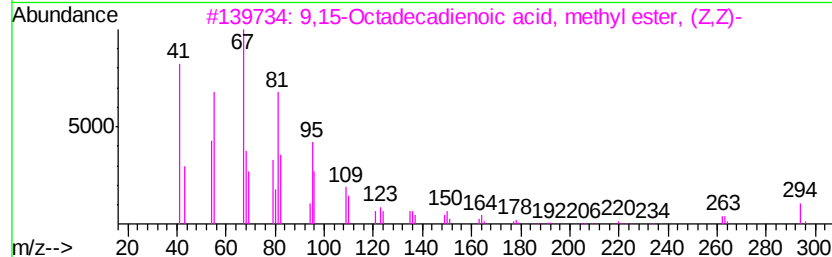
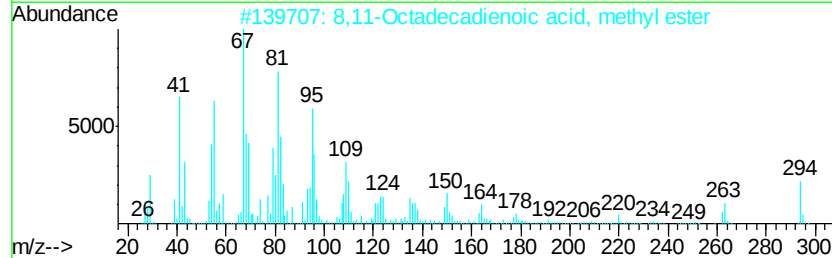
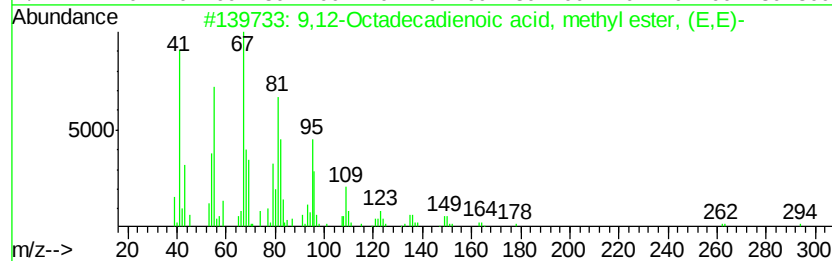
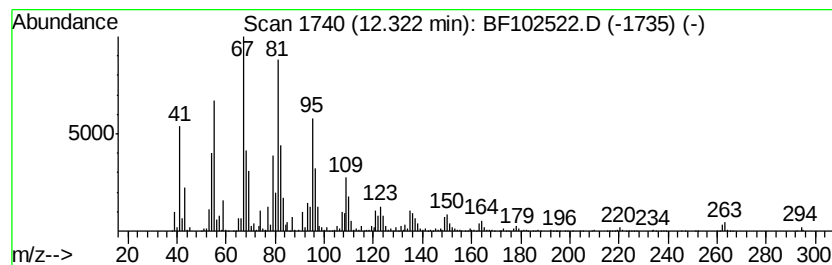
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	113.14 ng	5942310	Phenanthrene-d10	11.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2		8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	99
3		9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
4		10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
5		9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

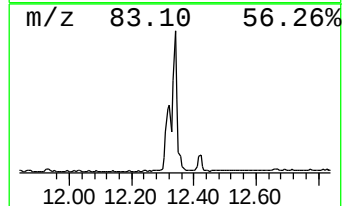
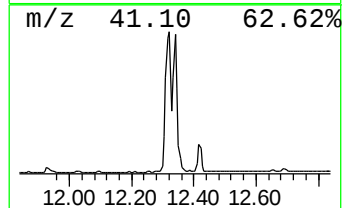
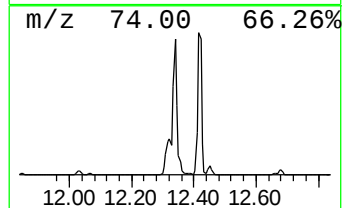
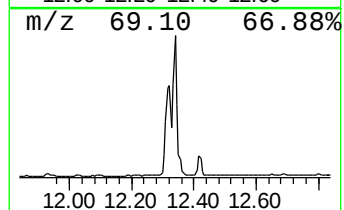
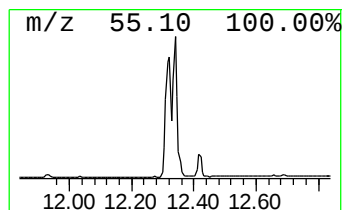
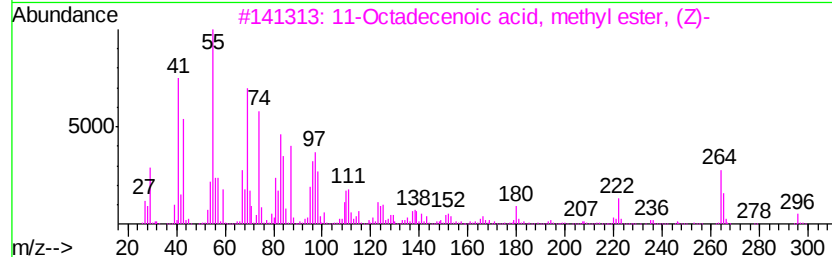
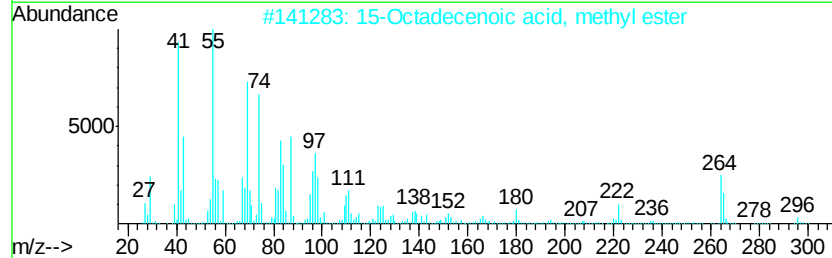
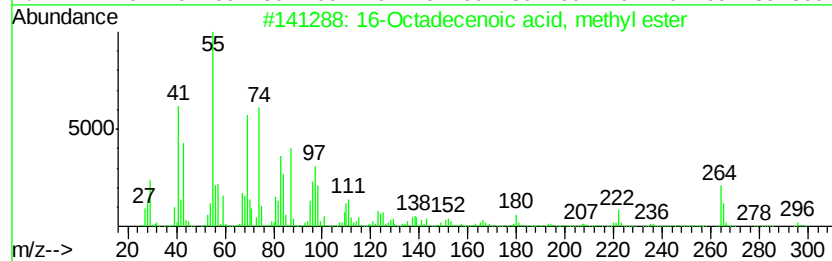
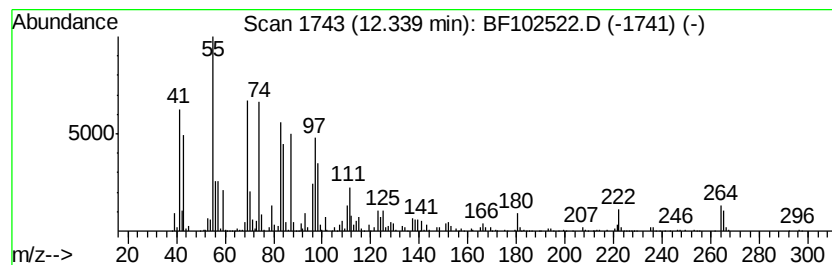
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 16-Octadecenoic acid, methyl ester Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.34	82.25 ng	4319720	Phenanthrene-d10	11.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	97
2		15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	93
3		11-Octadecenoic acid, methyl ester...	296	C19H36O2	001937-63-9	93
4		11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	87
5		12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

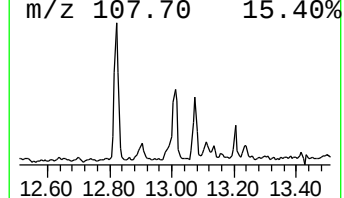
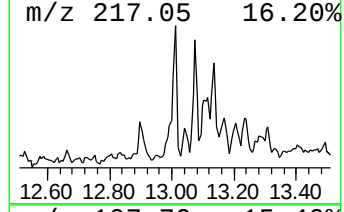
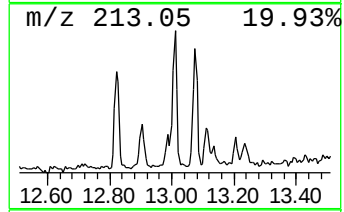
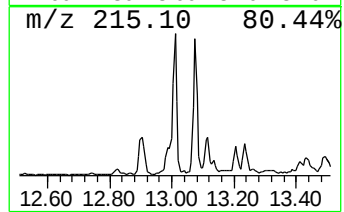
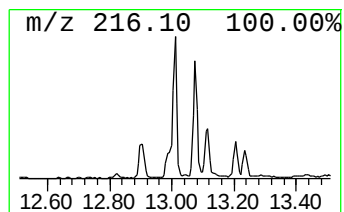
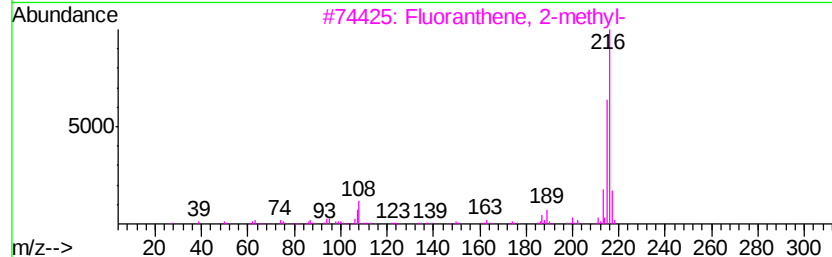
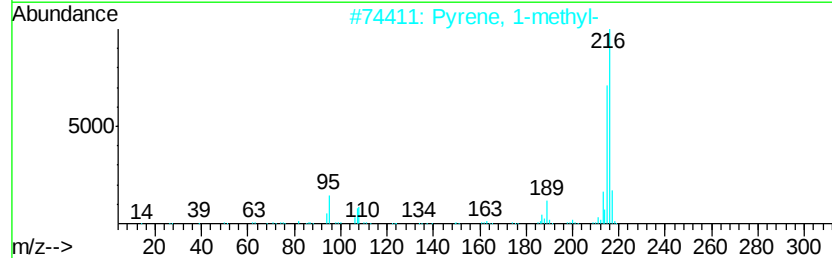
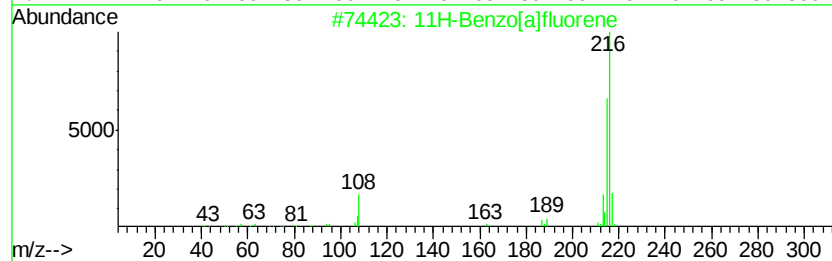
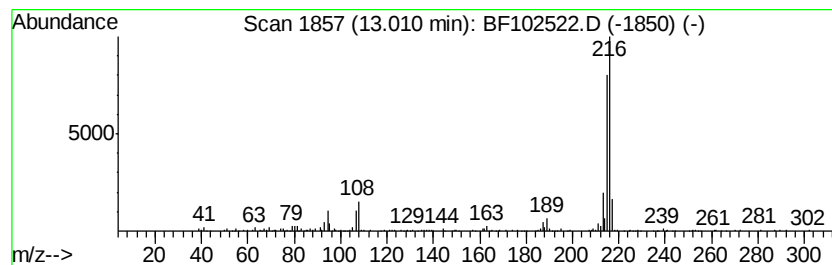
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 11H-Benzo[a]fluorene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.01	4.89 ng	311521	Chrysene-d12	13.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	94
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
3			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	93
4			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	81



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
Data File : BF102522.D
Acq On : 27 Jan 2018 20:42
Operator : SJ/JU
Sample : J1009-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-012518-B

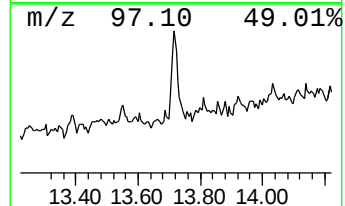
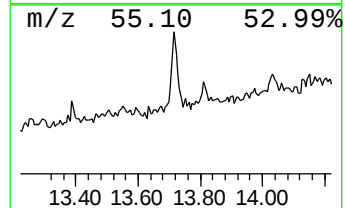
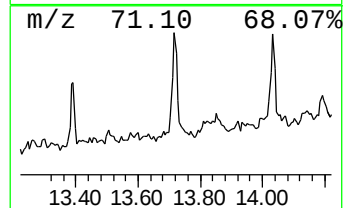
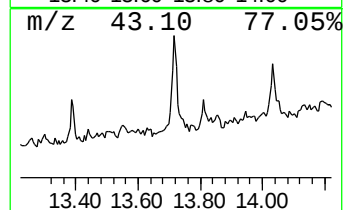
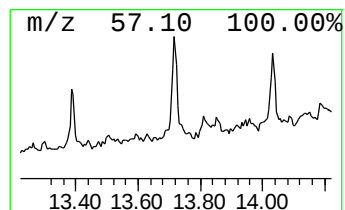
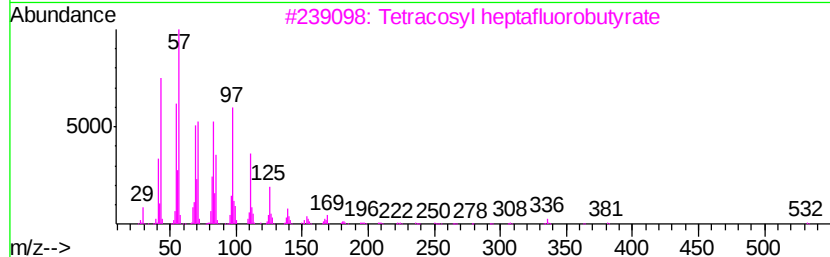
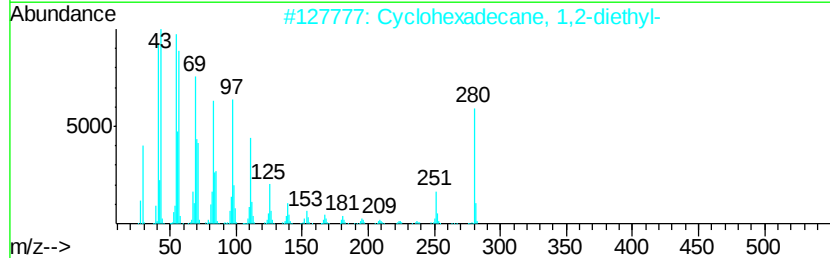
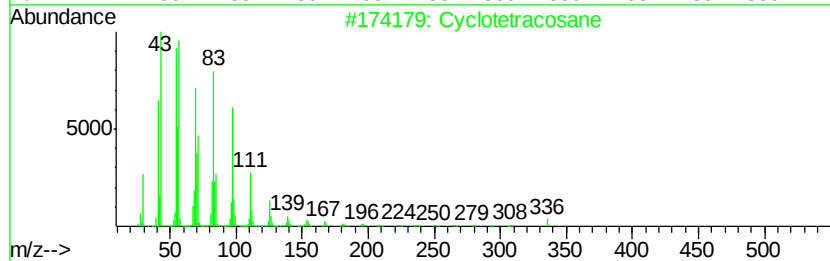
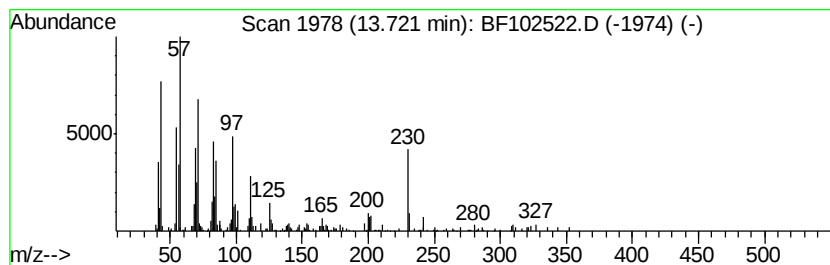
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Cyclotetracosane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	4.87 ng	310609	Chrysene-d12	13.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotetracosane	336	C24H48	000297-03-0	97
2		Cyclohexadecane, 1,2-diethyl-	280	C20H40	1000155-85-3	93
3		Tetracosyl heptafluorobutyrate	550	C28H49F7O2	1000351-83-7	91
4		Hexacosyl trifluoroacetate	478	C28H53F3O2	1000351-75-0	91
5		Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012718\
 Data File : BF102522.D
 Acq On : 27 Jan 2018 20:42
 Operator : SJ/JU
 Sample : J1009-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-06-012518-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.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
unknown6.49	6.49	73.7	ng	3719830	1	6.72	1009340	20.0
Undecane, 2,6-dim...	8.05	5.7	ng	365444	2	8.00	1286000	20.0
Cyclohexane, (4-m...	8.29	4.6	ng	295510	2	8.00	1286000	20.0
Nonane, 3-methyl-	8.42	9.4	ng	601902	2	8.00	1286000	20.0
Nonane, 5-butyl-	8.59	14.1	ng	907002	2	8.00	1286000	20.0
Dodecane, 2,6,10-...	9.02	7.5	ng	469408	3	9.75	1248580	20.0
Tetradecane	9.15	20.7	ng	1292240	3	9.75	1248580	20.0
Naphthalene, 2,7-...	9.41	4.3	ng	265575	3	9.75	1248580	20.0
Hexadecane	9.68	13.6	ng	850711	3	9.75	1248580	20.0
Pentadecane, 2,6,...	10.39	4.9	ng	306129	3	9.75	1248580	20.0
Heptadecane	10.65	15.8	ng	831119	4	11.24	1050400	20.0
Octadecane	11.10	5.8	ng	302968	4	11.24	1050400	20.0
Naphtho[2,3-b]thi...	11.13	4.2	ng	222302	4	11.24	1050400	20.0
Hexadecanoic acid...	11.63	42.2	ng	2214490	4	11.24	1050400	20.0
9,12-Octadecadien...	12.32	113.1	ng	5942310	4	11.24	1050400	20.0
16-Octadecenoic a...	12.34	82.3	ng	4319720	4	11.24	1050400	20.0
11H-Benzo[a]fluorene	13.01	4.9	ng	311521	5	13.88	1274870	20.0
Cyclotetracosane	13.72	4.9	ng	310609	5	13.88	1274870	20.0