

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF103017\
 Data File : BF099993.D
 Acq On : 31 Oct 2017 6:43
 Operator : SJ/JU
 Sample : I6201-02
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CF-1A(4-5)

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

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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.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	5.004	493	496	522	rBV	146207	266209	14.76%	1.407%
2	5.410	562	565	597	rBV	1517213	1723738	95.59%	9.112%
3	6.434	735	739	757	rBV	1584108	1756592	97.42%	9.285%
4	6.557	757	760	776	rVB	1872708	1793288	99.45%	9.479%
5	6.787	796	799	808	rBV	509322	451600	25.04%	2.387%
6	6.939	822	825	839	rVB	1528269	1371909	76.08%	7.252%
7	7.357	892	896	911	rBV	1130064	1058911	58.72%	5.597%
8	7.804	969	972	976	rBV3	23473	25437	1.41%	0.134%
9	7.934	989	994	996	rBV3	18516	20304	1.13%	0.107%
10	8.069	1014	1017	1023	rBV	717986	638380	35.40%	3.374%
11	8.116	1023	1025	1028	rVB2	45107	37586	2.08%	0.199%
12	8.151	1028	1031	1038	rVB6	14452	23022	1.28%	0.122%
13	8.263	1047	1050	1052	rVB3	26309	22229	1.23%	0.118%
14	8.351	1060	1065	1069	rVB5	30644	42776	2.37%	0.226%
15	8.475	1083	1086	1088	rBV	85241	69062	3.83%	0.365%
16	8.563	1099	1101	1104	rBV3	17865	22278	1.24%	0.118%
17	8.628	1109	1112	1115	rBV	39210	38577	2.14%	0.204%
18	8.745	1129	1132	1136	rVB2	26563	33112	1.84%	0.175%
19	8.833	1145	1147	1150	rBV4	16355	22551	1.25%	0.119%
20	8.886	1153	1156	1158	rBV2	61730	53878	2.99%	0.285%
21	8.963	1166	1169	1174	rVB4	23836	19892	1.10%	0.105%
22	9.075	1185	1188	1191	rBV	131900	108813	6.03%	0.575%
23	9.145	1195	1200	1203	rBV	1928882	1803194	100.00%	9.532%
24	9.210	1207	1211	1216	rBV4	65555	85340	4.73%	0.451%
25	9.251	1216	1218	1220	rVV2	23072	20231	1.12%	0.107%
26	9.286	1222	1224	1227	rVB4	39388	34554	1.92%	0.183%
27	9.328	1227	1231	1233	rBV5	23230	30498	1.69%	0.161%
28	9.416	1241	1246	1249	rBV6	26332	46954	2.60%	0.248%
29	9.480	1253	1257	1261	rVV6	24190	44333	2.46%	0.234%
30	9.539	1261	1267	1274	rVB2	356000	439315	24.36%	2.322%
31	9.598	1275	1277	1279	rBV	33914	30130	1.67%	0.159%
32	9.680	1288	1291	1296	rBV2	68054	79930	4.43%	0.423%
33	9.722	1296	1298	1301	rVB	43272	34803	1.93%	0.184%
34	9.822	1312	1315	1319	rBV	716952	664411	36.85%	3.512%

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35	9.939	1329	1335	1337	rBV2	27700	40925	2.27%	0.216%
36	10.004	1343	1346	1347	rBV2	27096	25285	1.40%	0.134%
37	10.069	1354	1357	1360	rVB2	61317	59292	3.29%	0.313%
38	10.139	1365	1369	1373	rBV	82153	96466	5.35%	0.510%
39	10.186	1373	1377	1379	rVB4	27919	32126	1.78%	0.170%
40	10.222	1380	1383	1386	rBV2	58234	59557	3.30%	0.315%
41	10.251	1386	1388	1390	rVV	32016	21826	1.21%	0.115%
42	10.369	1402	1408	1409	rBV4	43360	62402	3.46%	0.330%
43	10.439	1418	1420	1421	rBV	39332	29284	1.62%	0.155%
44	10.463	1421	1424	1426	rVB	132533	115764	6.42%	0.612%
45	10.516	1431	1433	1435	rBV3	23614	25419	1.41%	0.134%
46	10.539	1435	1437	1441	rVB	118555	90253	5.01%	0.477%
47	10.622	1447	1451	1464	rBV	1168263	1158447	64.24%	6.124%
48	10.727	1465	1469	1477	rVB	301912	301644	16.73%	1.594%
49	10.810	1481	1483	1486	rBV2	24777	24112	1.34%	0.127%
50	10.957	1505	1508	1511	rVB2	40313	36960	2.05%	0.195%
51	11.039	1520	1522	1524	rBV3	19918	23512	1.30%	0.124%
52	11.192	1544	1548	1551	rBV	136238	134129	7.44%	0.709%
53	11.322	1566	1570	1574	rBV	768137	706684	39.19%	3.736%
54	11.545	1605	1608	1611	rBV3	44140	40616	2.25%	0.215%
55	11.833	1653	1657	1666	rBV4	57146	96827	5.37%	0.512%
56	12.374	1746	1749	1753	rVB4	20102	20781	1.15%	0.110%
57	12.710	1803	1806	1813	rVB2	37683	34448	1.91%	0.182%
58	12.904	1835	1839	1842	rBV	1620542	1464303	81.21%	7.740%
59	13.416	1923	1926	1930	rVB	22362	20194	1.12%	0.107%
60	13.786	1985	1989	1998	rBV	101237	126732	7.03%	0.670%
61	13.980	2018	2022	2029	rBV	501182	497529	27.59%	2.630%
62	14.115	2041	2045	2049	rBV	19928	21074	1.17%	0.111%
63	14.427	2094	2098	2101	rBV	26307	26158	1.45%	0.138%
64	14.739	2148	2151	2157	rVB	30541	29140	1.62%	0.154%
65	15.068	2203	2207	2212	rVB	23921	28112	1.56%	0.149%
66	15.439	2266	2270	2272	rBV	24445	29030	1.61%	0.153%
67	15.474	2272	2276	2289	rVB	314561	416472	23.10%	2.201%
68	15.862	2337	2342	2347	rVB2	21517	28097	1.56%	0.149%
69	16.351	2421	2425	2433	rVB6	16771	31047	1.72%	0.164%
70	16.921	2517	2522	2528	rBV4	10347	19876	1.10%	0.105%
71	17.603	2632	2638	2644	rBV8	12397	29521	1.64%	0.156%

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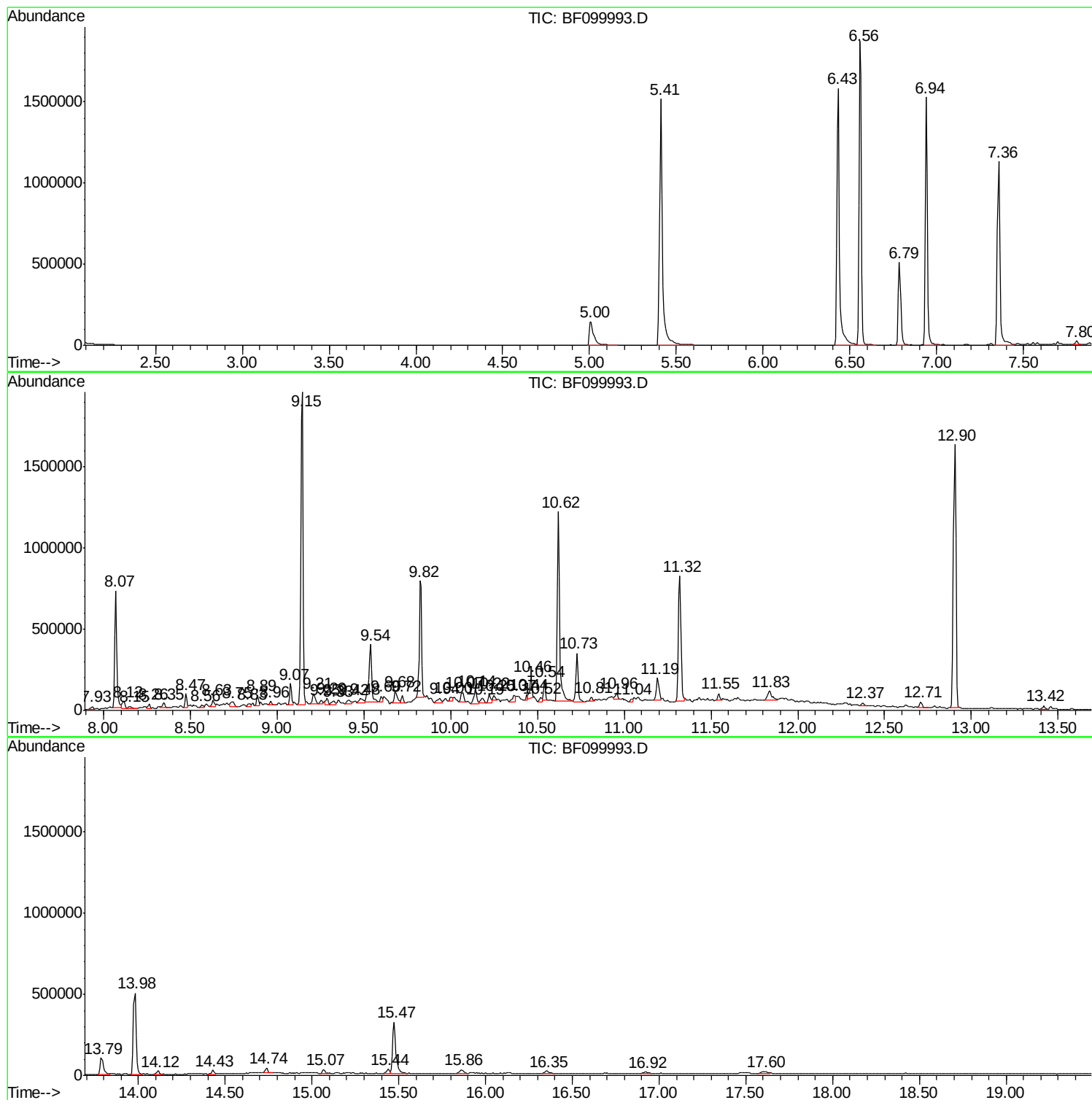
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Misc :
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Instrument :
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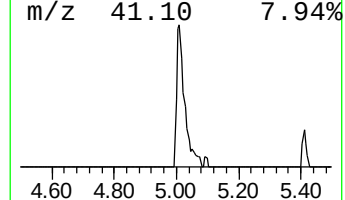
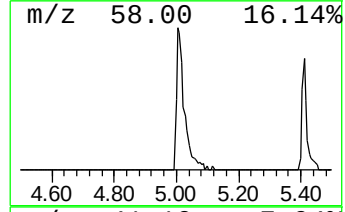
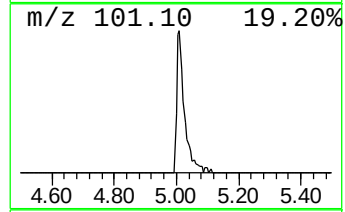
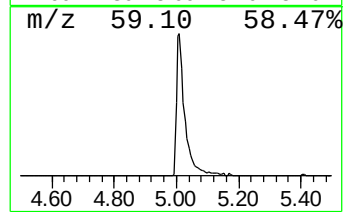
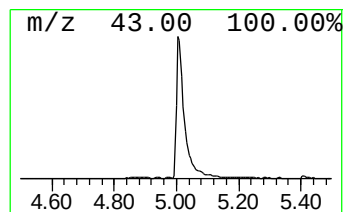
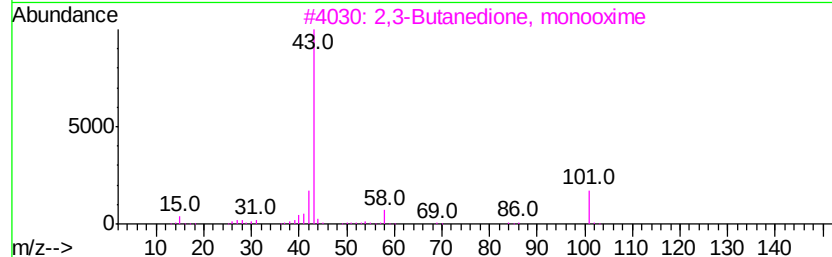
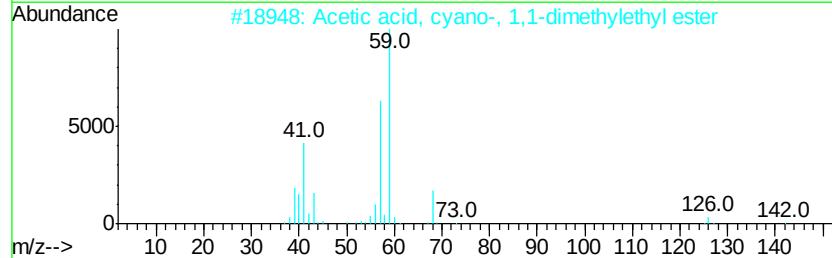
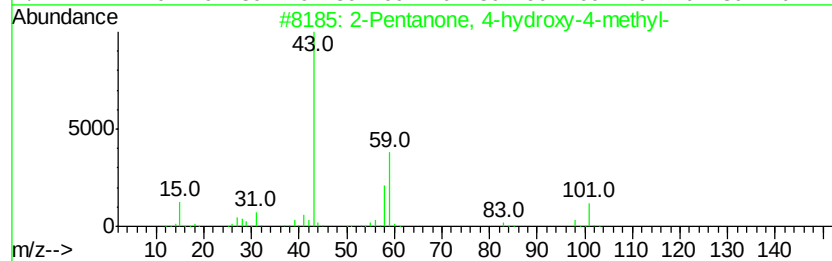
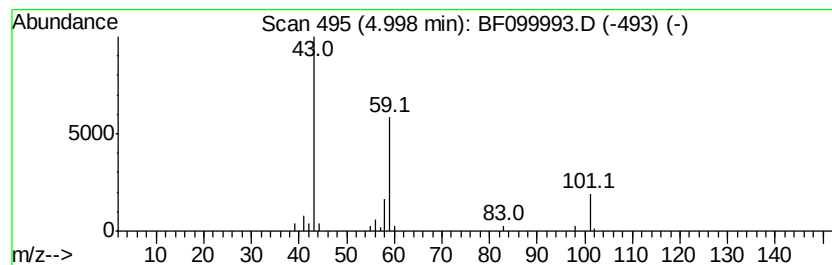
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	11.79 ng	266209	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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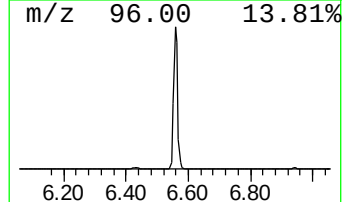
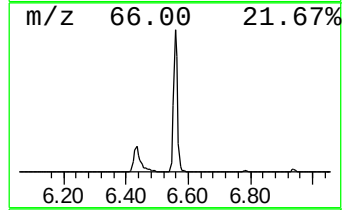
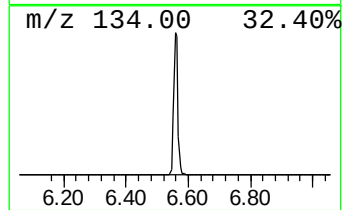
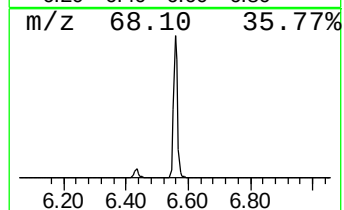
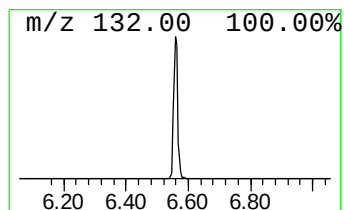
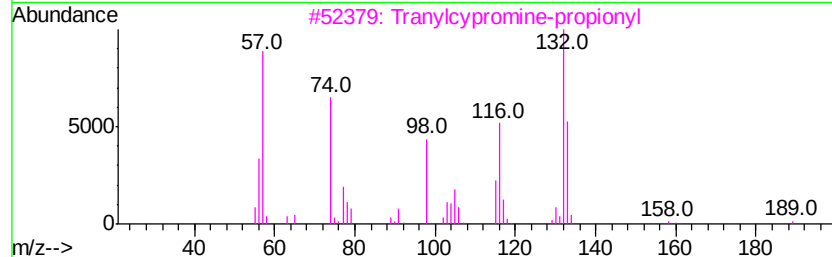
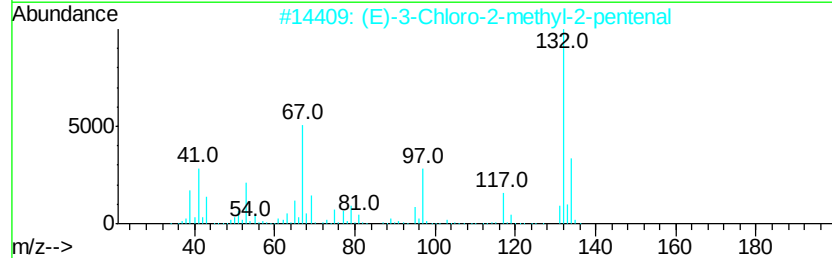
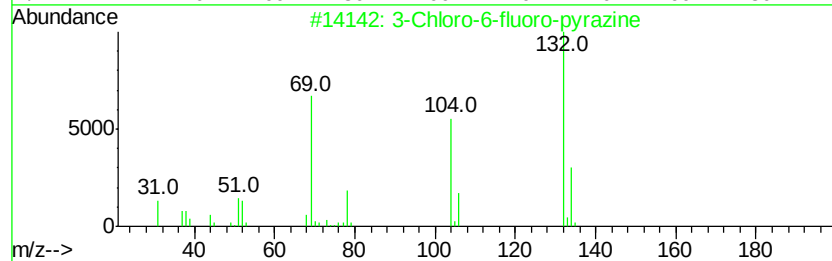
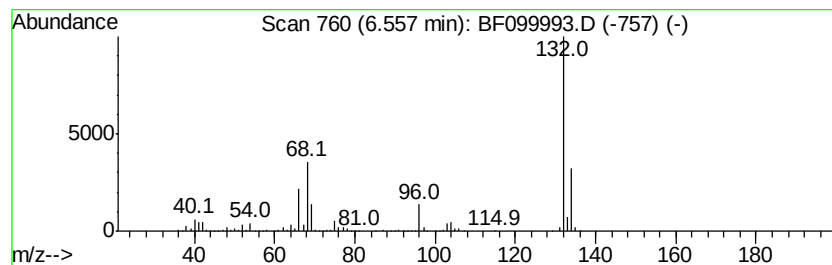
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	79.42 ng	1793290	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9



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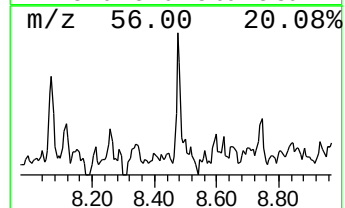
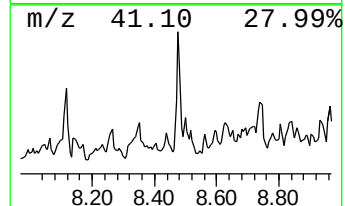
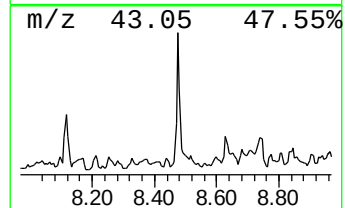
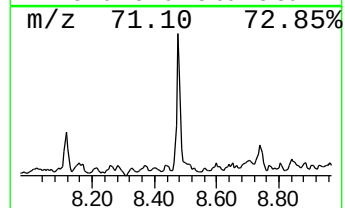
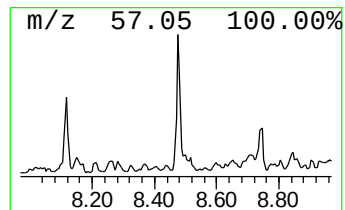
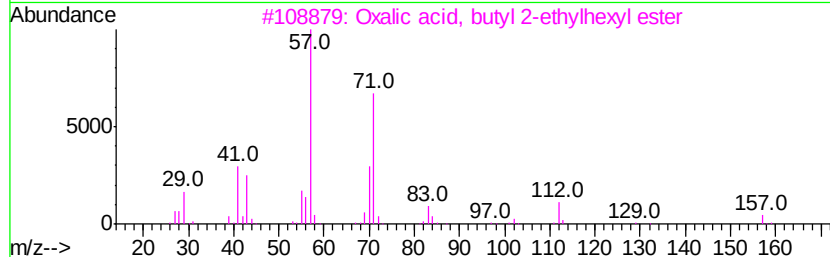
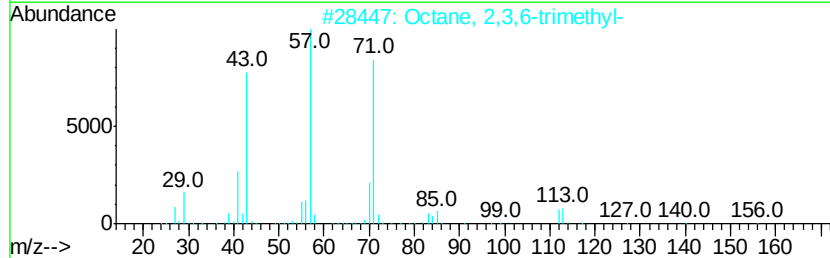
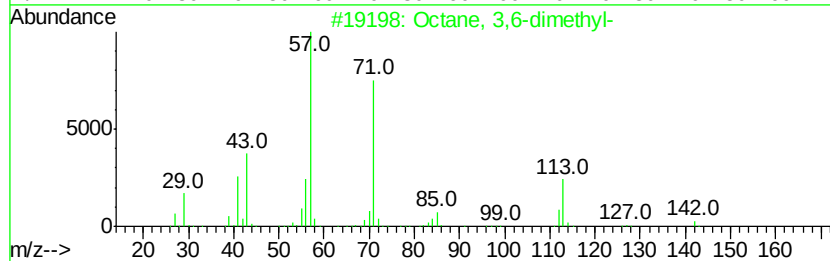
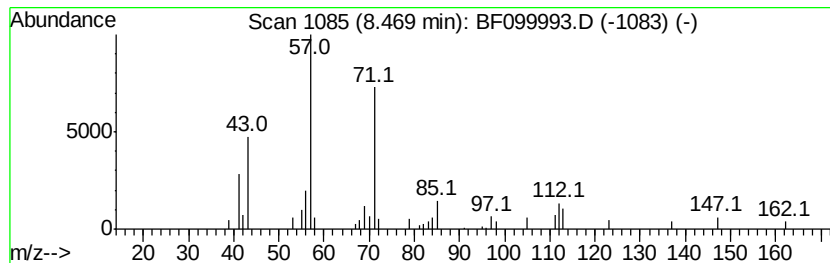
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Octane, 3,6-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.47	2.16 ng	69062	Naphthalene-d8	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	72
2		Octane, 2,3,6-trimethyl-	156	C11H24	062016-33-5	72
3		Oxalic acid, butyl 2-ethylhexyl ...	258	C14H26O4	1000309-38-6	64
4		Pentadecane, 7-methyl-	226	C16H34	006165-40-8	64
5		Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	64



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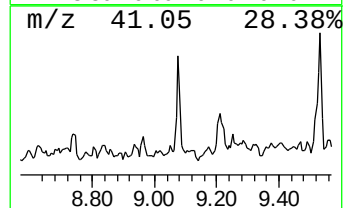
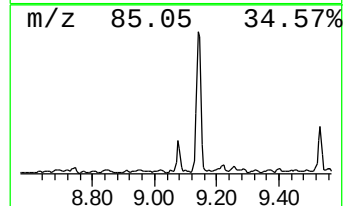
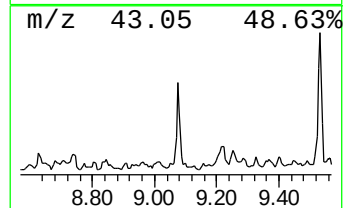
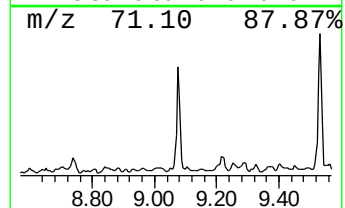
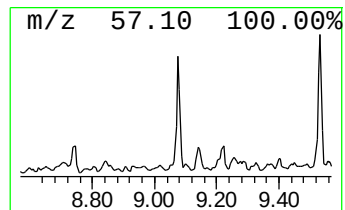
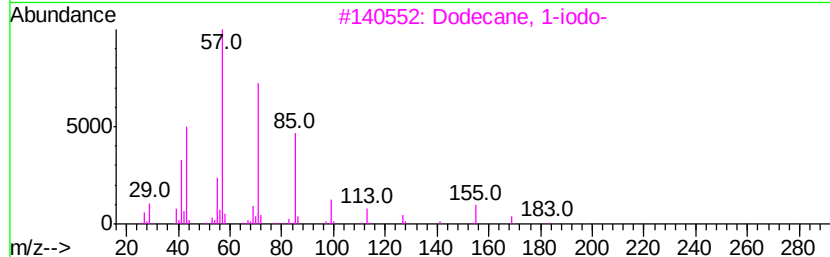
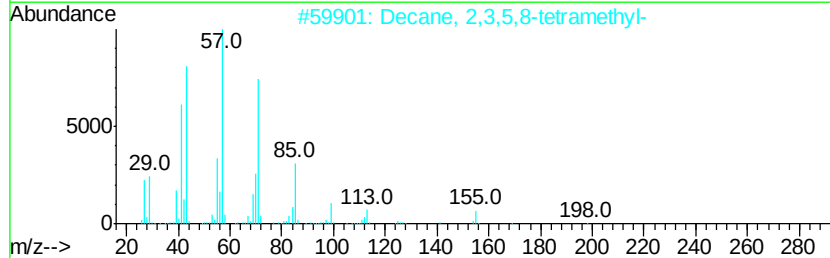
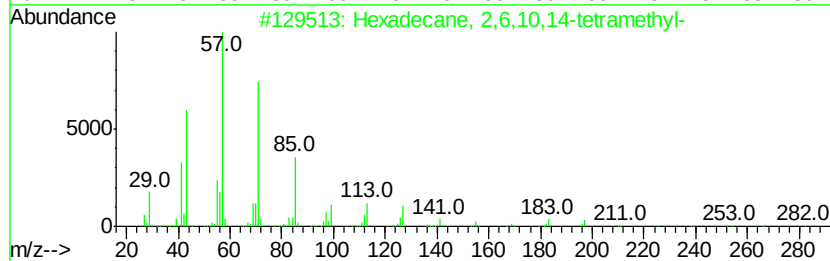
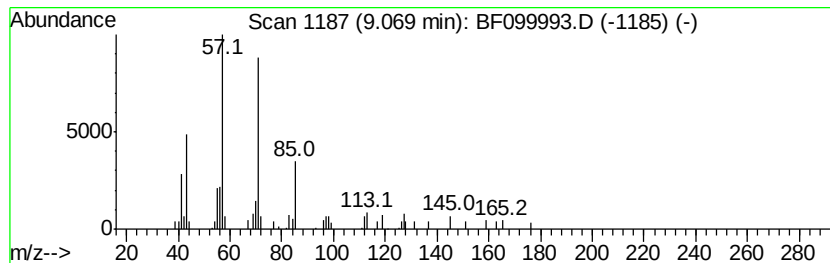
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Hexadecane, 2,6,10,14-tetra... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.07	3.28 ng	108813	Acenaphthene-d10	9.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	78
2		Decane, 2,3,5,8-tetramethyl-	198	C14H30	192823-15-7	72
3		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	64
4		Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	64
5		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF103017\
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Operator : SJ/JU
Sample : I6201-02
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ALS Vial : 42 Sample Multiplier: 1

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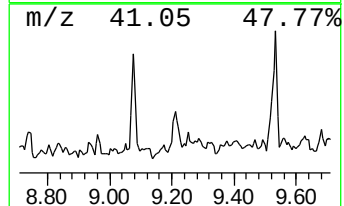
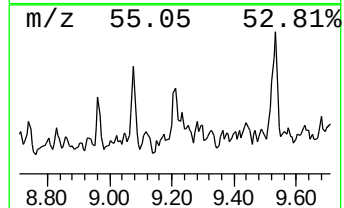
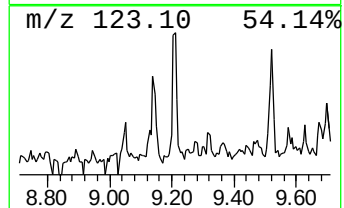
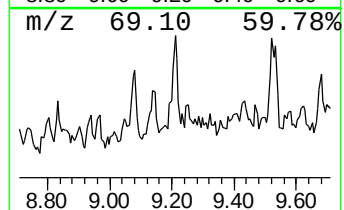
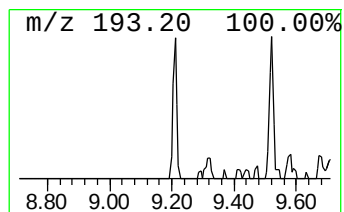
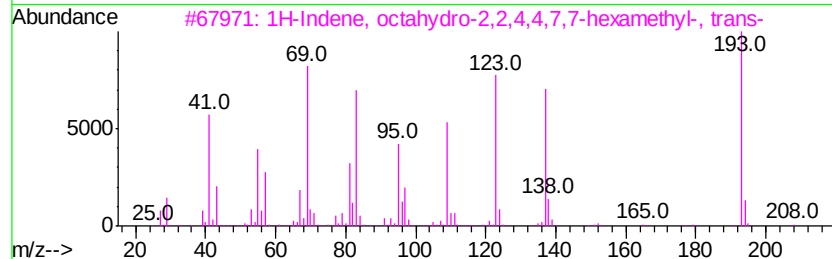
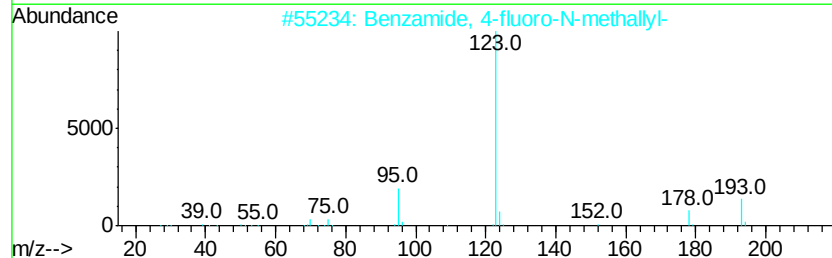
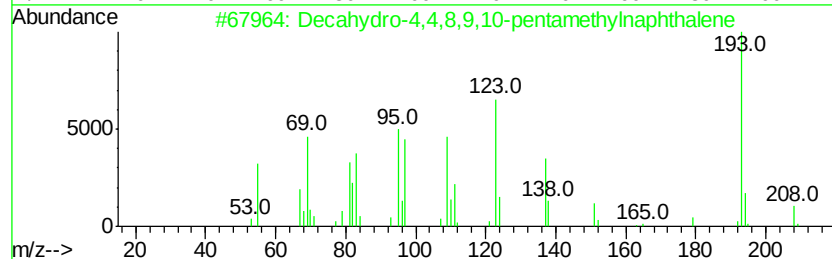
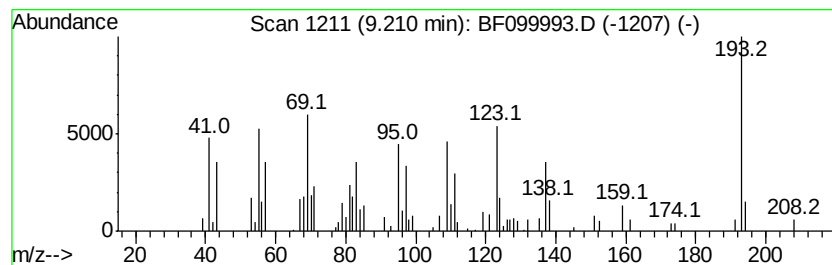
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Decahydro-4,4,8,9,10-pentam... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.21	2.57 ng	85340	Acenaphthene-d10	9.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	64
2		Benzamide, 4-fluoro-N-methyl-	193	C11H12FNO	1000340-31-9	47
3		1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	43
4		1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	43
5		2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	38



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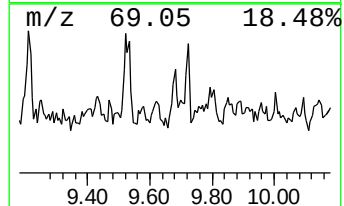
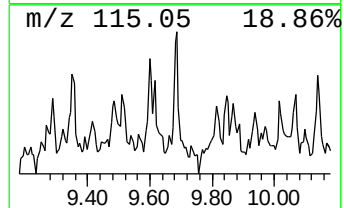
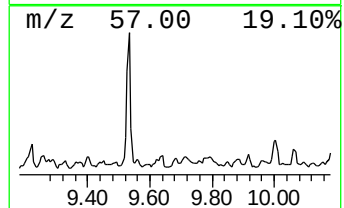
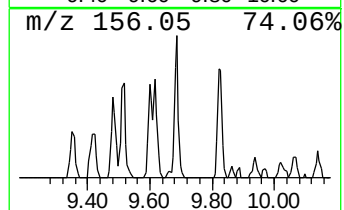
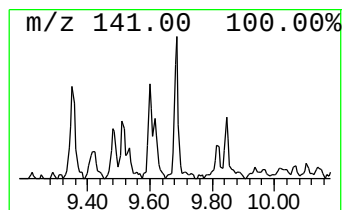
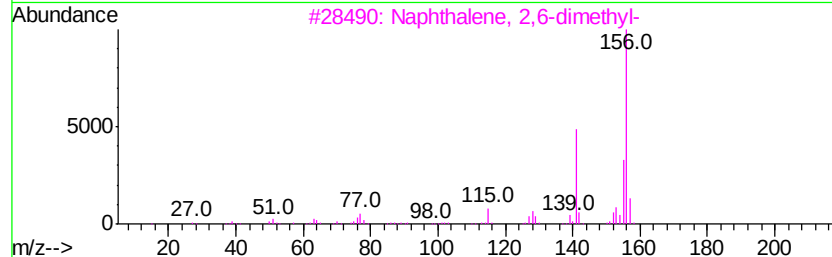
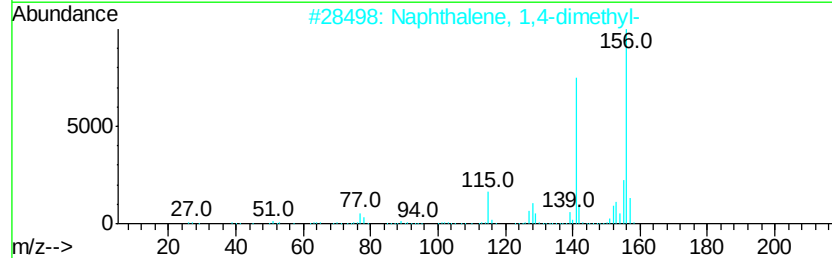
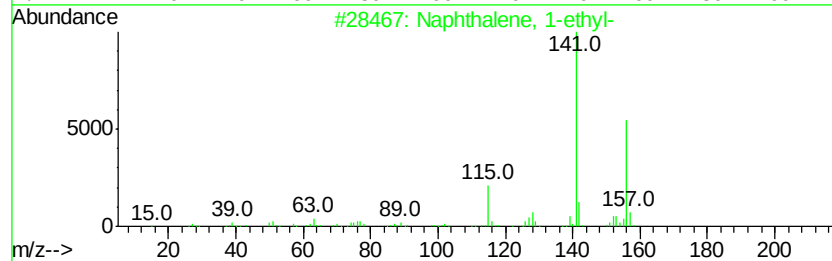
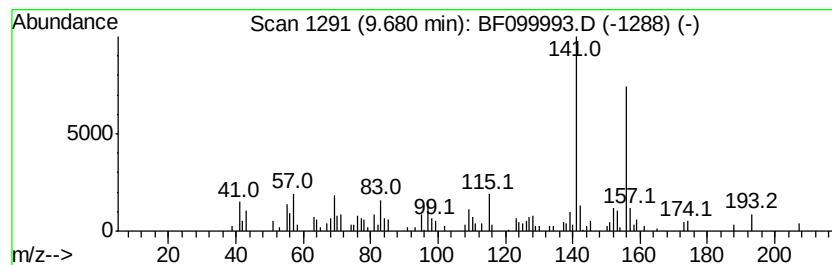
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Naphthalene, 1-ethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.68	2.41 ng	79930	Acenaphthene-d10	9.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	87
2		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	87
3		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	87
4		Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	83
5		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	83



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Data File : BF099993.D
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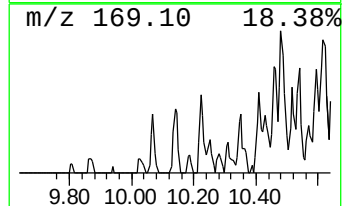
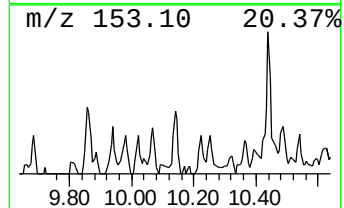
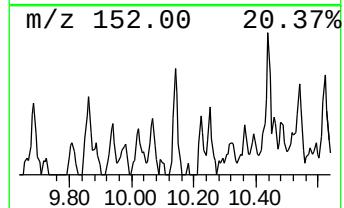
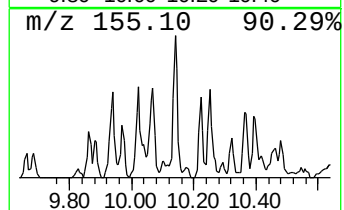
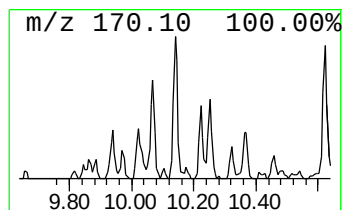
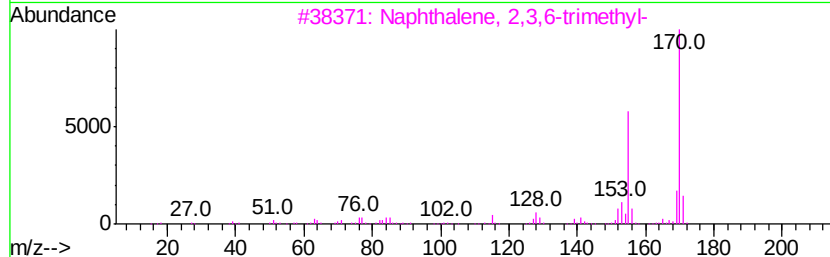
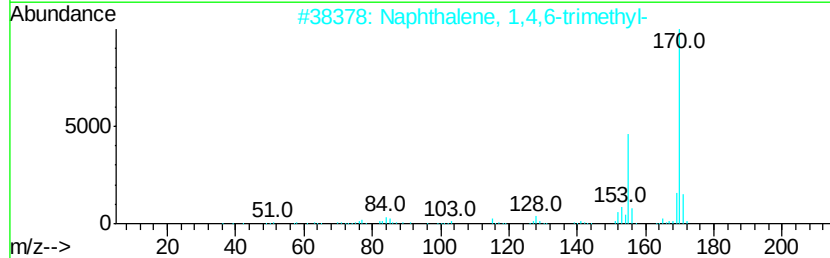
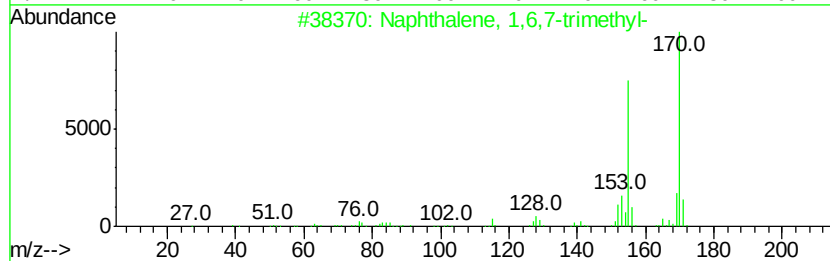
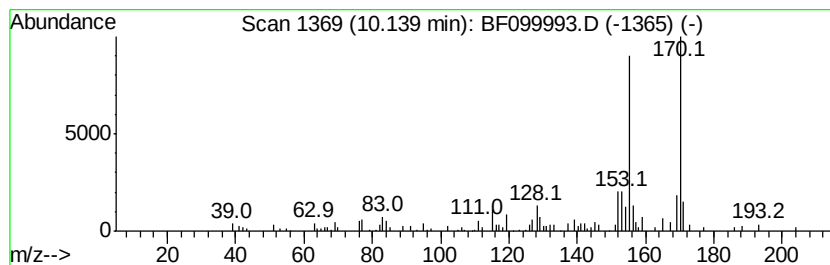
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphthalene, 1,6,7-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.14	2.90 ng	96466	Acenaphthene-d10	9.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
2		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	94
3		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	94
4		4,6,8-Trimethylazulene	170	C13H14	000941-81-1	87
5		3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF103017\
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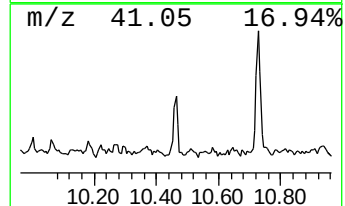
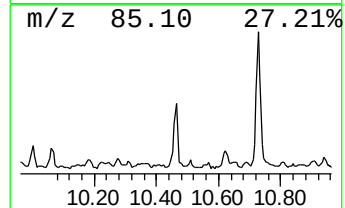
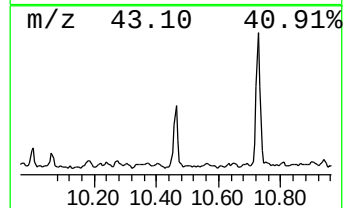
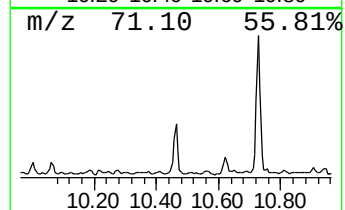
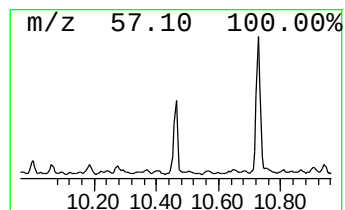
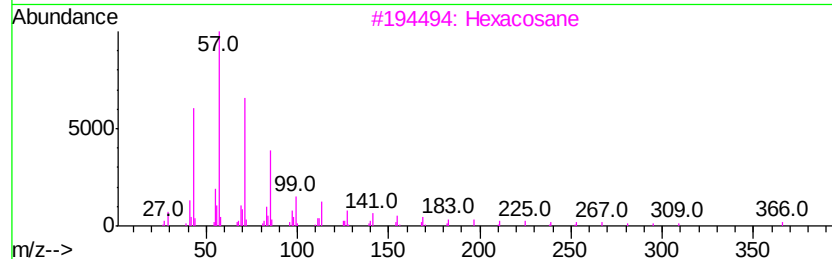
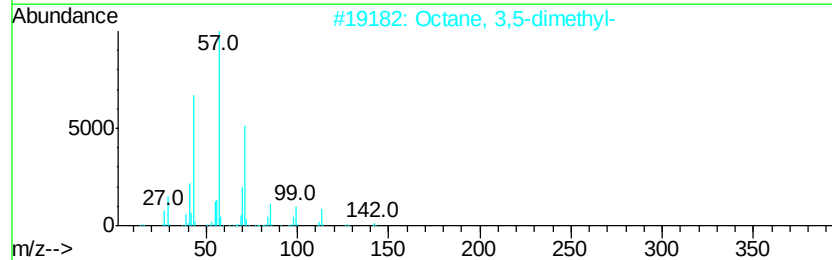
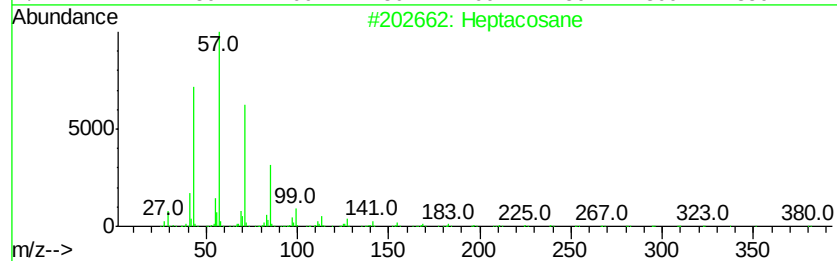
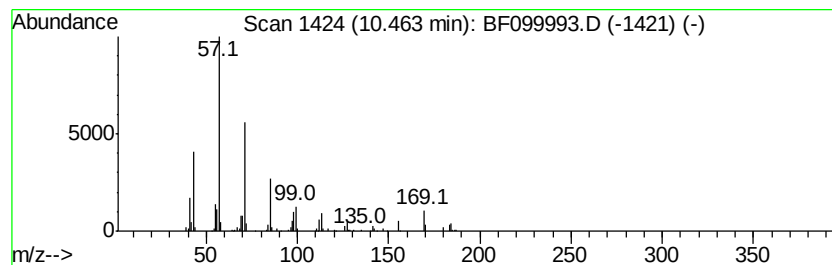
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.46	3.48 ng	115764	Acenaphthene-d10	9.82
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptacosane	380 C27H56	000593-49-7	72
2	Octane, 3,5-dimethyl-	142 C10H22	015869-93-9	68
3	Hexacosane	366 C26H54	000630-01-3	64
4	Dodecane, 1-iodo-	296 C12H25I	004292-19-7	58
5	Octane, 3,6-dimethyl-	142 C10H22	015869-94-0	55



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF103017\
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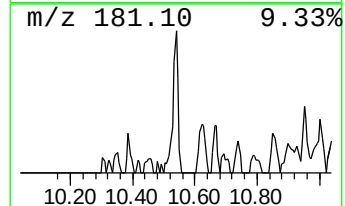
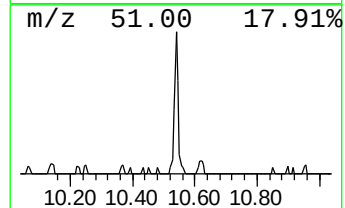
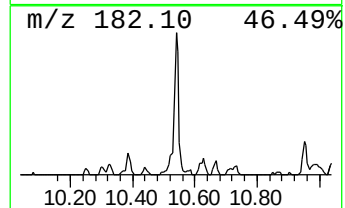
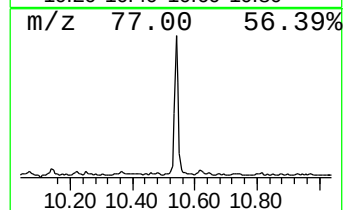
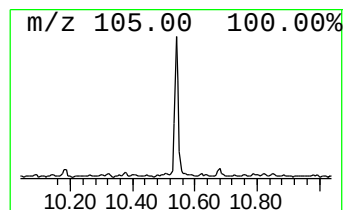
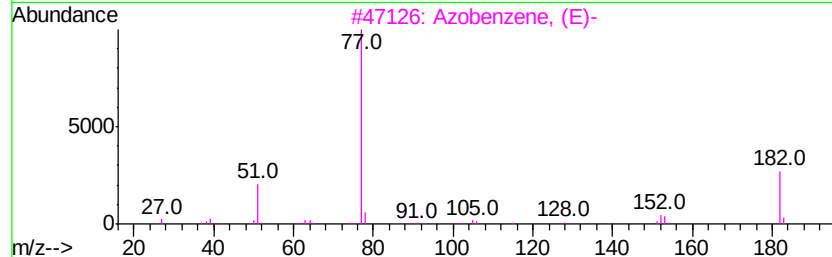
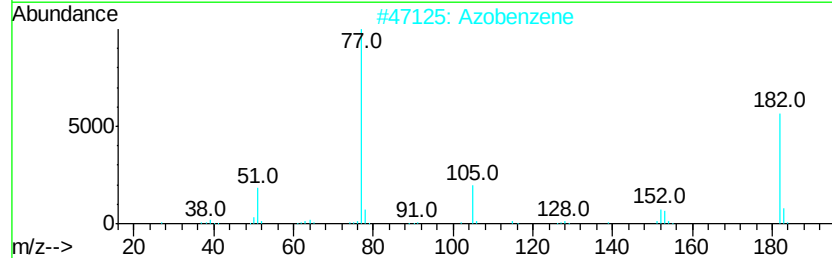
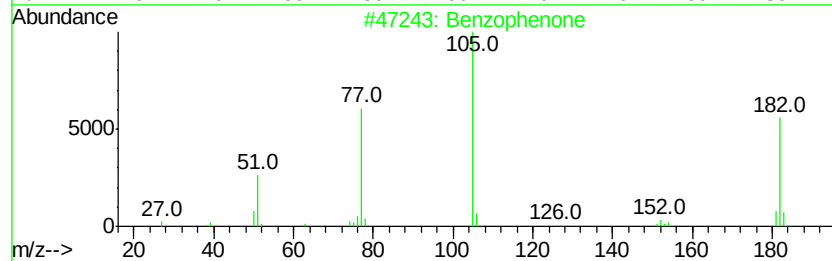
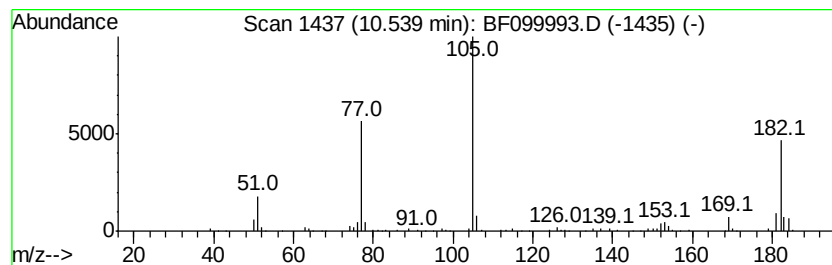
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzophenone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.54	2.72 ng	90253	Acenaphthene-d10		9.82	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzophenone	182	C13H10O	000119-61-9	95
2		Azobenzene	182	C12H10N2	000103-33-3	76
3		Azobenzene, (E)-	182	C12H10N2	017082-12-1	53
4		2-(2-Oxo-2-phenyl-ethyl)-malonon...	184	C11H8N2O	1000296-76-9	38
5		Phenyl 4-pyridyl ketone	183	C12H9NO	014548-46-0	35



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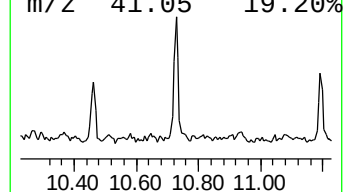
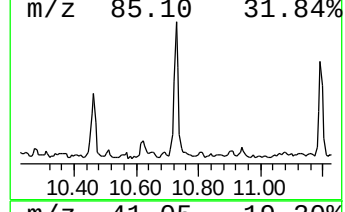
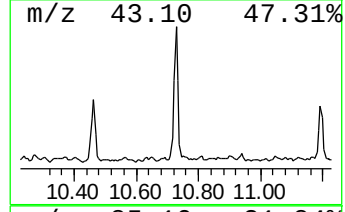
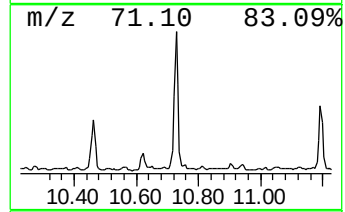
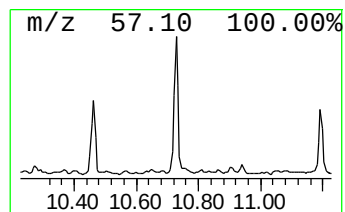
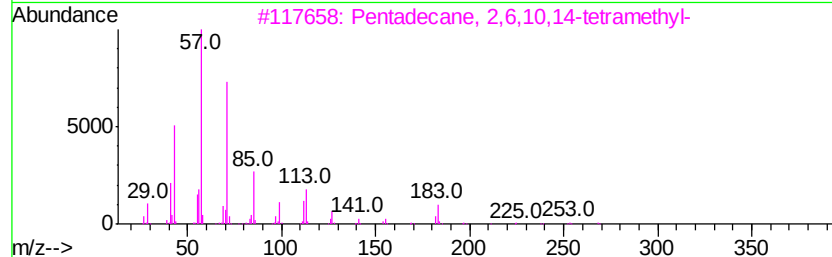
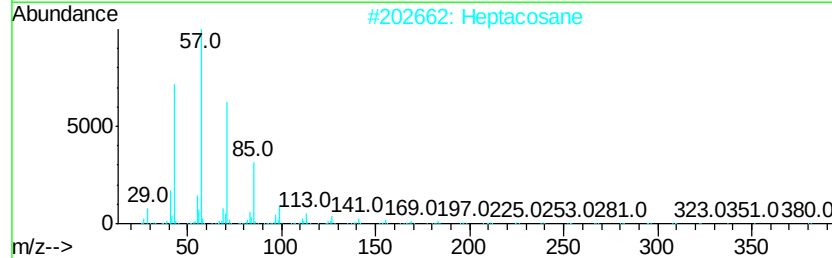
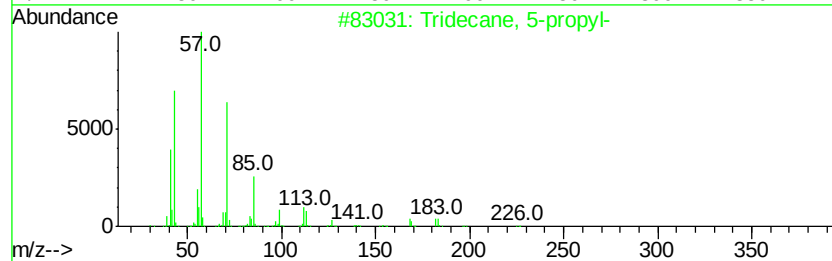
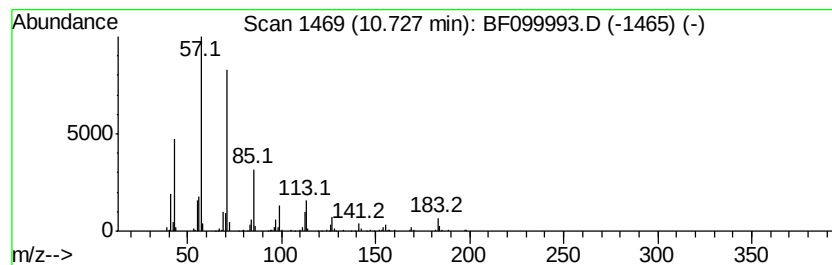
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tridecane, 5-propyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.73	8.54 ng	301644	Phenanthrene-d10		11.32
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 5-propyl-	226	C16H34	055045-11-9	90
2	Heptacosane	380	C27H56	000593-49-7	86
3	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	83
4	Hexadecane, 2,6,10-trimethyl-	268	C19H40	055000-52-7	80
5	Dodecane	170	C12H26	000112-40-3	72



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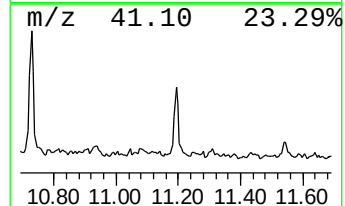
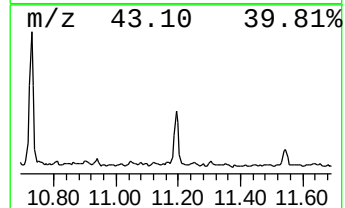
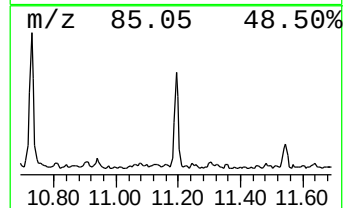
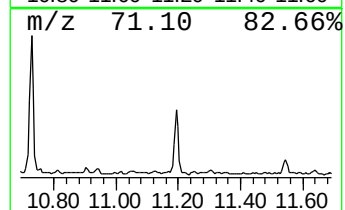
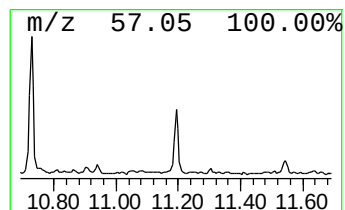
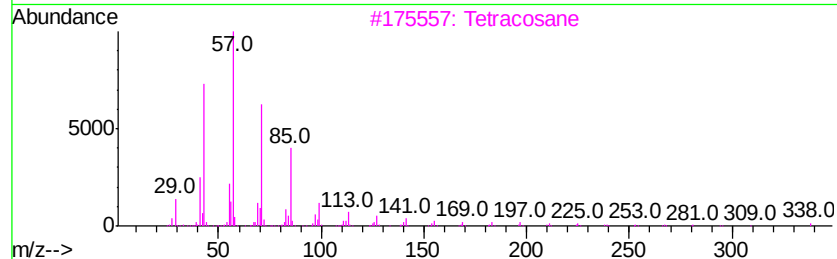
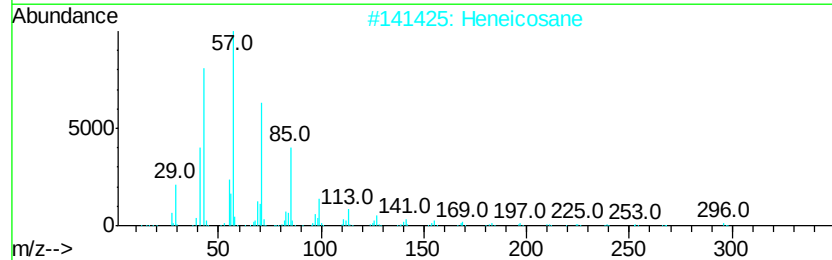
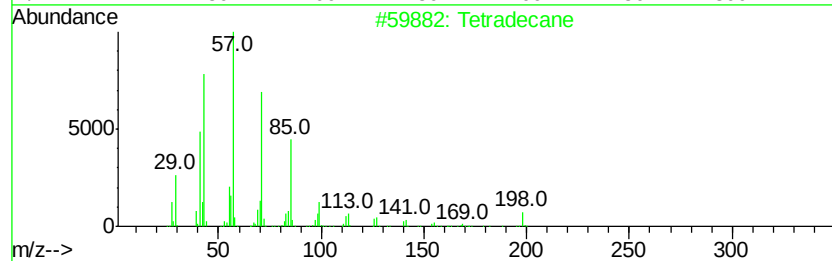
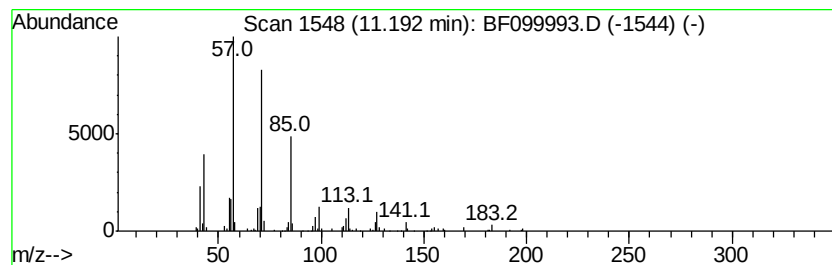
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tetradecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.19	3.80 ng	134129	Phenanthrene-d10	11.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	90
2			Heneicosane	296	C21H44	000629-94-7	90
3			Tetracosane	338	C24H50	000646-31-1	90
4			Heptadecane	240	C17H36	000629-78-7	86
5			Octadecane	254	C18H38	000593-45-3	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF103017\
Data File : BF099993.D
Acq On : 31 Oct 2017 6:43
Operator : SJ/JU
Sample : I6201-02
Misc :
ALS Vial : 42 Sample Multiplier: 1

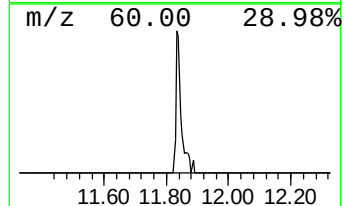
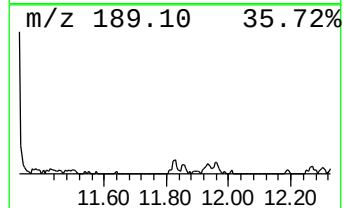
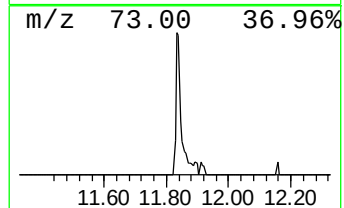
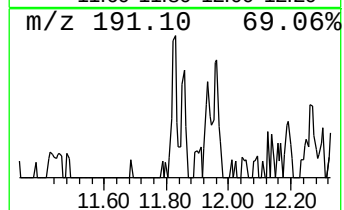
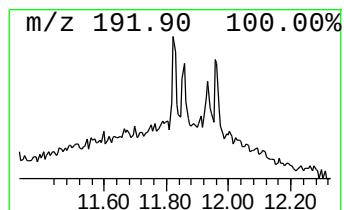
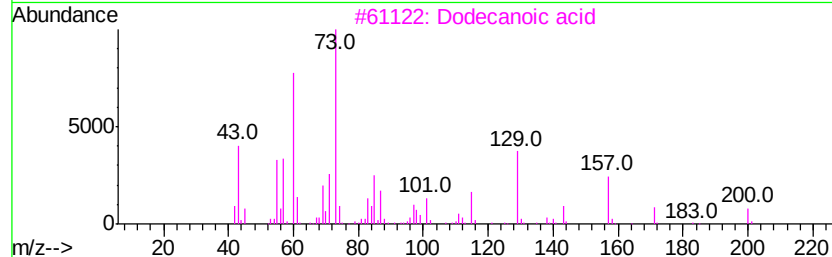
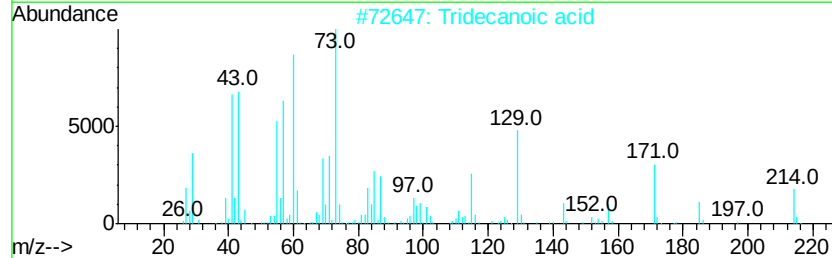
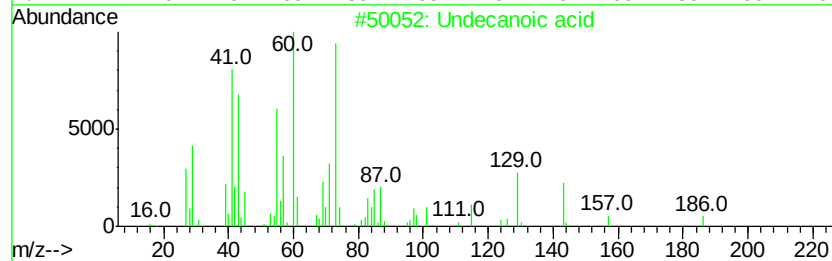
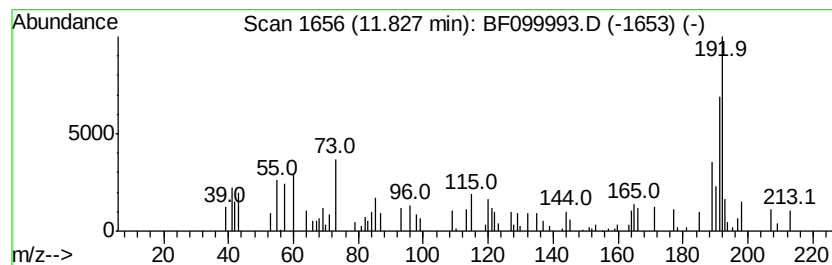
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ClientSampleId :
CF-1A(4-5)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Undecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.83	2.74 ng	96827	Phenanthrene-d10		11.32
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecanoic acid	186	C11H22O2	000112-37-8	59
2	Tridecanoic acid	214	C13H26O2	000638-53-9	53
3	Dodecanoic acid	200	C12H24O2	000143-07-7	53
4	Heptanoic acid, propyl ester	172	C10H20O2	007778-87-2	50
5	n-Decanoic acid	172	C10H20O2	000334-48-5	50



Data Path : Z:\HPCHEM1\BNA F\DATA\BF103017\
Data File : BF099993.D
Acq On : 31 Oct 2017 6:43
Operator : SJ/JU
Sample : I6201-02
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CF-1A(4-5)

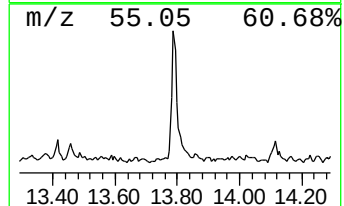
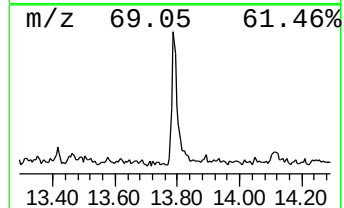
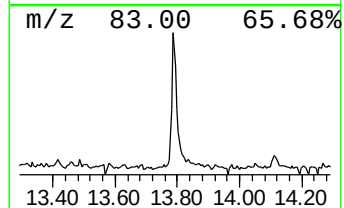
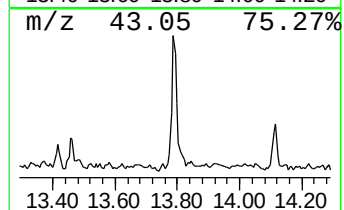
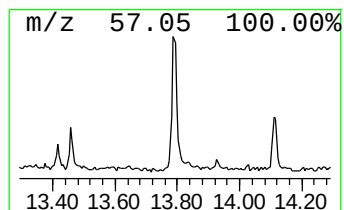
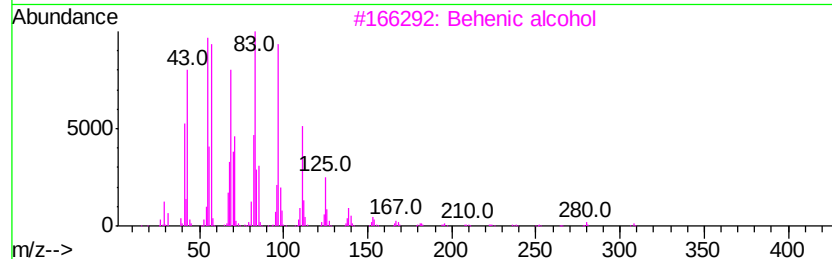
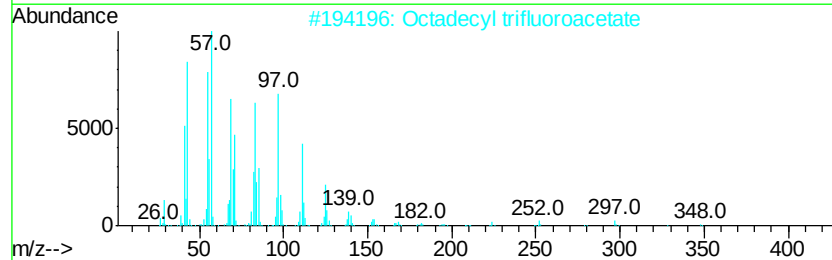
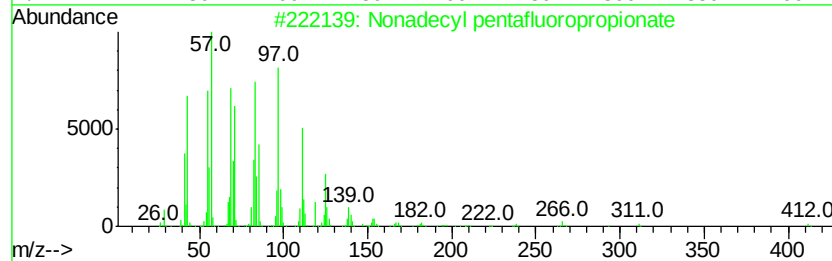
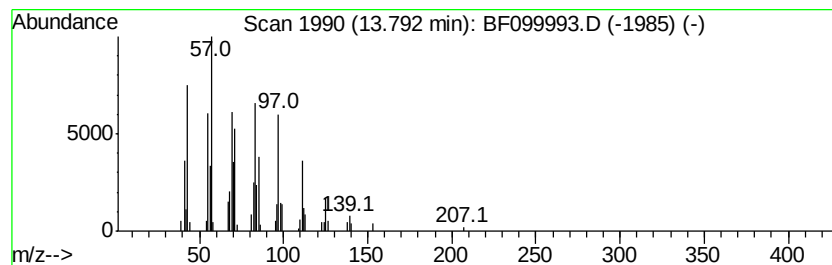
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Nonadecyl pentafluoropropio... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	5.09 ng	126732	Chrysene-d12	13.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
2		Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
3		Behenic alcohol	326	C22H46O	000661-19-8	87
4		Tricosyl heptafluorobutyrte	536	C27H47F7O2	1000351-83-4	87
5		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF103017\
 Data File : BF099993.D
 Acq On : 31 Oct 2017 6:43
 Operator : SJ/JU
 Sample : I6201-02
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CF-1A(4-5)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.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
2-Pentanone, 4-hy...	5.00	11.8	ng	266209	1	6.79	451600	20.0
unknown6.56	6.56	79.4	ng	1793290	1	6.79	451600	20.0
Octane, 3,6-dimet...	8.47	2.2	ng	69062	2	8.07	638380	20.0
Hexadecane, 2,6,1...	9.07	3.3	ng	108813	3	9.82	664411	20.0
Decahydro-4,4,8,9...	9.21	2.6	ng	85340	3	9.82	664411	20.0
Naphthalene, 1-et...	9.68	2.4	ng	79930	3	9.82	664411	20.0
Naphthalene, 1,6,...	10.14	2.9	ng	96466	3	9.82	664411	20.0
Heptacosane	10.46	3.5	ng	115764	3	9.82	664411	20.0
Benzophenone	10.54	2.7	ng	90253	3	9.82	664411	20.0
Tridecane, 5-propyl-	10.73	8.5	ng	301644	4	11.32	706684	20.0
Tetradecane	11.19	3.8	ng	134129	4	11.32	706684	20.0
Undecanoic acid	11.83	2.7	ng	96827	4	11.32	706684	20.0
Nonadecyl pentafl...	13.79	5.1	ng	126732	5	13.98	497529	20.0