

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF110117\
 Data File : BF100058.D
 Acq On : 1 Nov 2017 21:37
 Operator : SJ/JU
 Sample : I6236-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 133

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	4.345	381	384	410	rBV	253169	375785	17.81%	2.029%
2	4.992	489	494	510	rBV	866533	1112937	52.75%	6.008%
3	5.404	561	564	586	rBV	543569	758201	35.94%	4.093%
4	6.428	734	738	756	rBV	1337209	1446155	68.55%	7.807%
5	6.551	756	759	766	rVV	1190222	1048541	49.70%	5.660%
6	6.781	795	798	807	rBV	474349	470141	22.28%	2.538%
7	6.934	820	824	839	rBV	1622647	1437292	68.13%	7.759%
8	7.351	892	895	912	rBV	1115871	1056355	50.07%	5.702%
9	8.063	1013	1016	1029	rVB	766395	659879	31.28%	3.562%
10	8.628	1109	1112	1113	rBV	31269	28546	1.35%	0.154%
11	9.139	1195	1199	1202	rBV	2405024	2109724	100.00%	11.389%
12	9.539	1264	1267	1273	rVB	276208	237769	11.27%	1.284%
13	9.739	1299	1301	1304	rVB	44748	30291	1.44%	0.164%
14	9.822	1312	1315	1319	rBV2	777837	640220	30.35%	3.456%
15	9.857	1319	1321	1324	rVB	56687	44056	2.09%	0.238%
16	10.033	1347	1351	1354	rBV2	37888	36366	1.72%	0.196%
17	10.239	1383	1386	1389	rVB	154785	124492	5.90%	0.672%
18	10.374	1406	1409	1417	rVB	55973	61376	2.91%	0.331%
19	10.463	1418	1424	1426	rBV	65557	79955	3.79%	0.432%
20	10.539	1434	1437	1442	rVB	124166	101889	4.83%	0.550%
21	10.627	1448	1452	1461	rBV2	153612	187778	8.90%	1.014%
22	10.733	1465	1470	1476	rBV2	123541	215891	10.23%	1.165%
23	10.869	1490	1493	1495	rBV2	23627	24129	1.14%	0.130%
24	10.916	1499	1501	1503	rBV2	22885	23038	1.09%	0.124%
25	11.180	1544	1546	1548	rBV	80655	71694	3.40%	0.387%
26	11.216	1549	1552	1554	rVV	116834	124726	5.91%	0.673%
27	11.239	1554	1556	1559	rVB2	34726	30845	1.46%	0.167%
28	11.345	1570	1574	1576	rBV	619155	561396	26.61%	3.031%
29	11.369	1576	1578	1583	rVV	338497	312460	14.81%	1.687%
30	11.427	1586	1588	1593	rVB	87819	86792	4.11%	0.469%
31	11.586	1613	1615	1617	rBV	38631	28171	1.34%	0.152%
32	11.639	1622	1624	1630	rVB	73690	65713	3.11%	0.355%
33	11.751	1641	1643	1648	rVB2	35195	31498	1.49%	0.170%
34	11.892	1664	1667	1669	rBV	40604	43103	2.04%	0.233%

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.921	1670	1672	1678	rVB2	43872	52293	2.48%	0.282%
36	11.980	1678	1682	1685	rBV5	28071	46747	2.22%	0.252%
37	12.016	1686	1688	1691	rVB	46253	43206	2.05%	0.233%
38	12.098	1699	1702	1705	rBV	66360	66374	3.15%	0.358%
39	12.545	1775	1778	1781	rBV2	92265	109580	5.19%	0.592%
40	12.716	1804	1807	1813	rVB	343497	346159	16.41%	1.869%
41	12.904	1835	1839	1842	rBV5	29934	39458	1.87%	0.213%
42	12.998	1848	1855	1862	rBV	271193	410079	19.44%	2.214%
43	13.180	1880	1886	1889	rBV	1388288	1590010	75.37%	8.583%
44	13.421	1925	1927	1930	rVB	40819	39951	1.89%	0.216%
45	13.480	1934	1937	1939	rBV	44815	42797	2.03%	0.231%
46	13.510	1939	1942	1945	rVB	31384	30231	1.43%	0.163%
47	13.957	2015	2018	2021	rBV	38890	36685	1.74%	0.198%
48	14.433	2094	2099	2106	rBV5	81573	150767	7.15%	0.814%
49	14.633	2127	2133	2137	rBV2	445123	670365	31.78%	3.619%
50	14.674	2137	2140	2145	rVB	107286	137234	6.50%	0.741%
51	14.786	2157	2159	2163	rVB2	25668	33777	1.60%	0.182%
52	14.921	2179	2182	2185	rBV2	31003	34771	1.65%	0.188%
53	15.415	2264	2266	2273	rBV4	37135	51660	2.45%	0.279%
54	16.239	2402	2406	2410	rBV	99529	164587	7.80%	0.888%
55	16.421	2434	2437	2441	rVB2	50089	64047	3.04%	0.346%
56	16.645	2473	2475	2480	rVB	57738	77498	3.67%	0.418%
57	16.733	2487	2490	2495	rVB	82764	99634	4.72%	0.538%
58	16.809	2499	2503	2509	rVB2	276294	394333	18.69%	2.129%
59	17.586	2632	2635	2643	rVB3	75891	125002	5.93%	0.675%

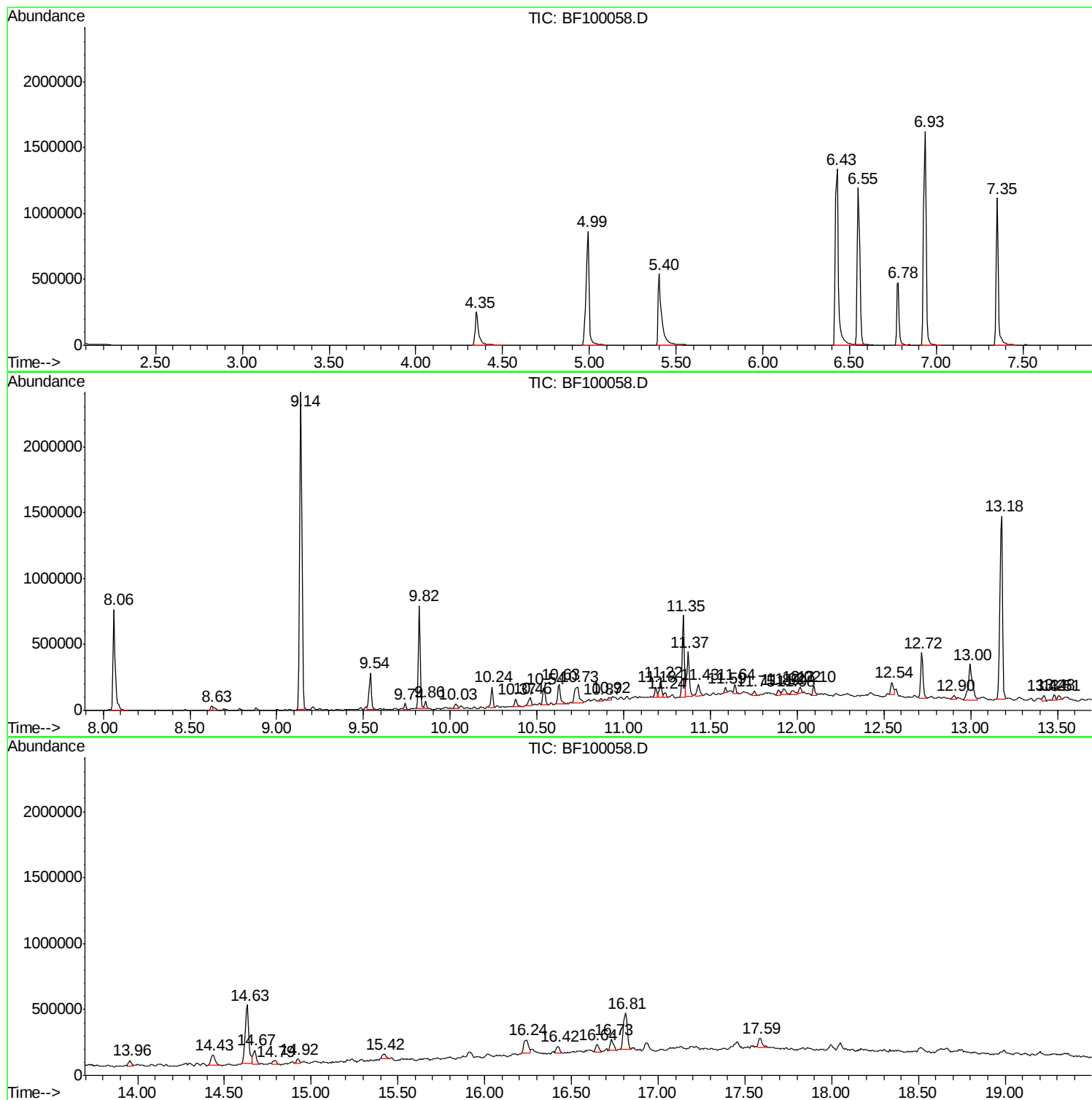
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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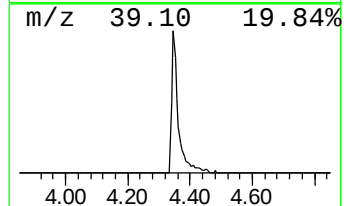
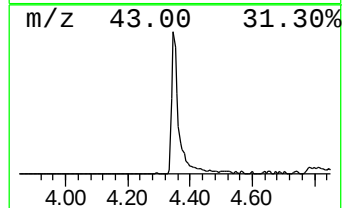
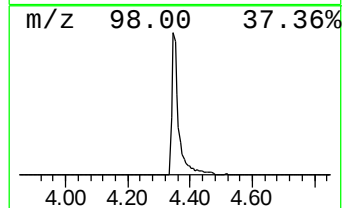
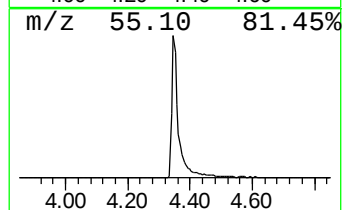
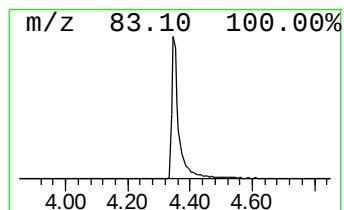
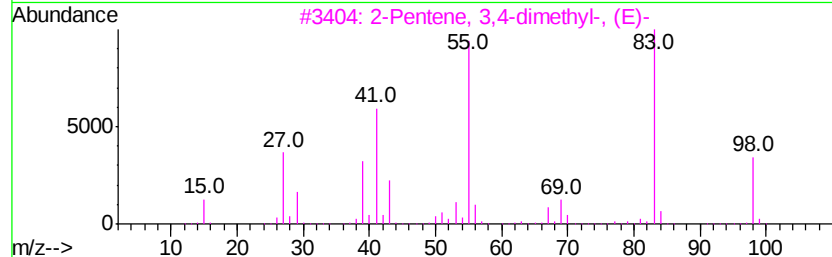
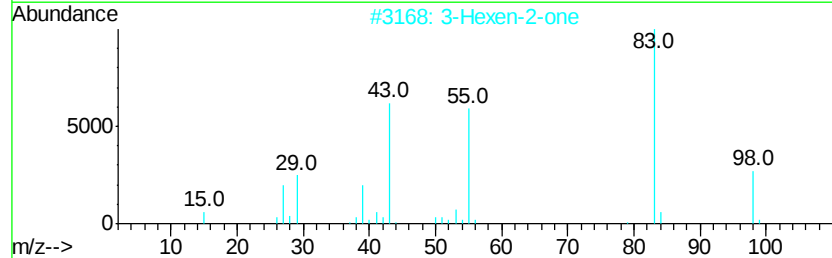
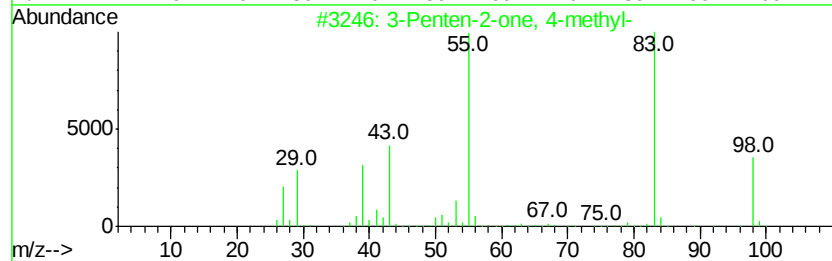
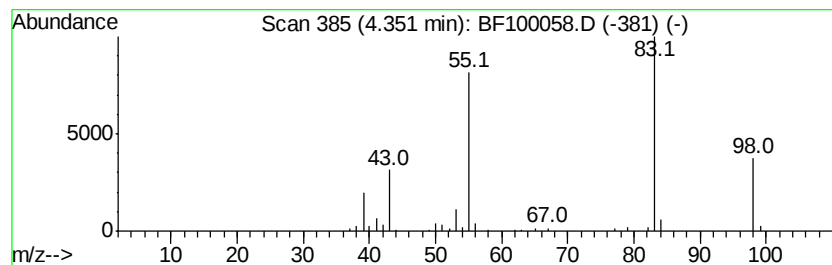
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.35	15.99 ng	375785	1,4-Dichlorobenzene-d4	6.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2		3-Hexen-2-one	98	C6H10O	000763-93-9	91
3		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	90
5		2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	86



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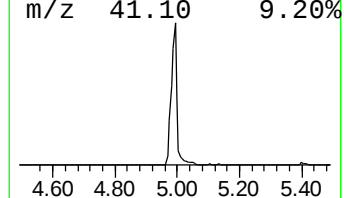
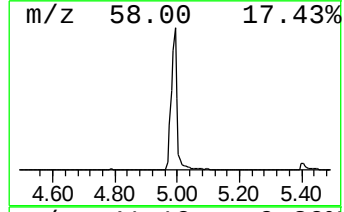
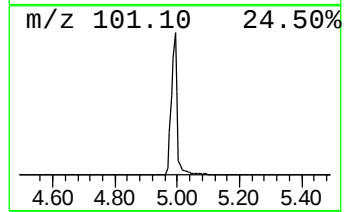
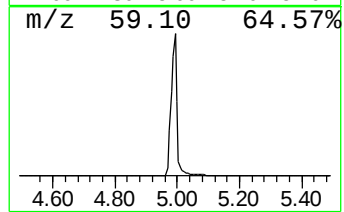
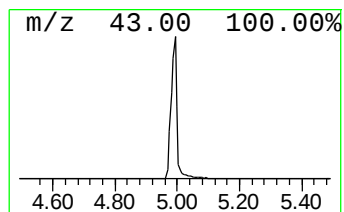
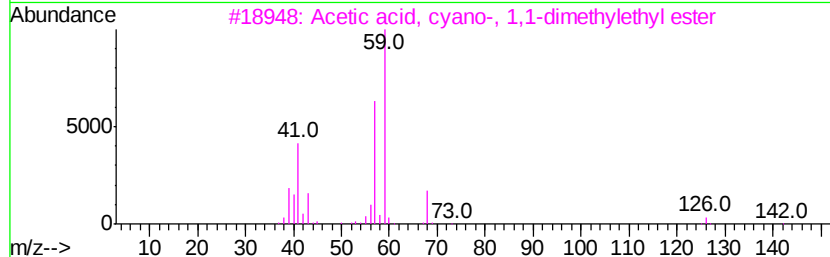
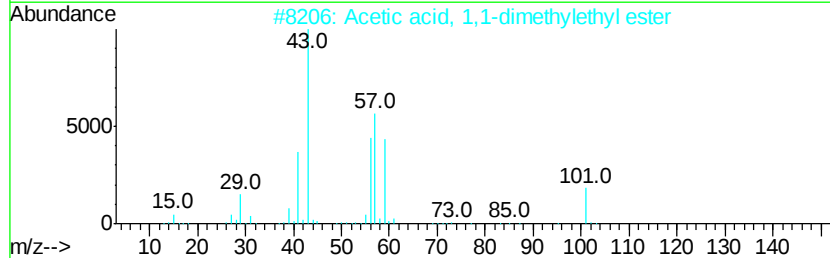
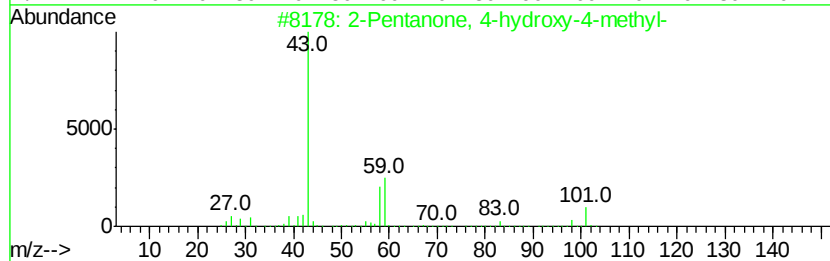
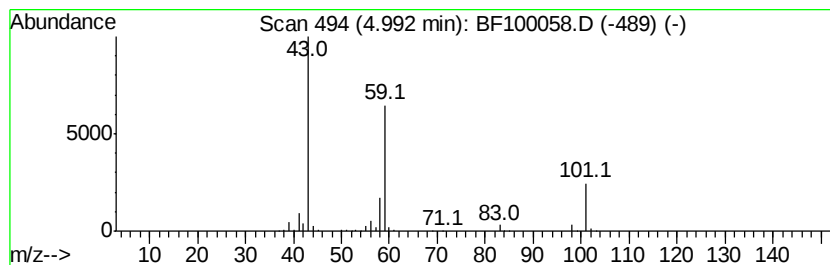
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	47.34 ng	1112940	1,4-Dichlorobenzene-d4	6.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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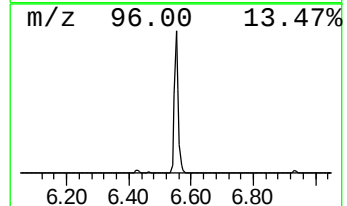
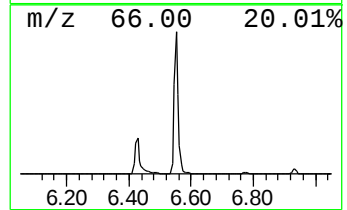
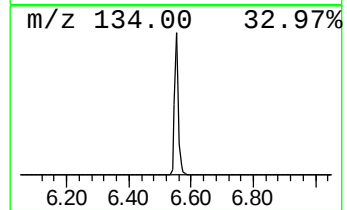
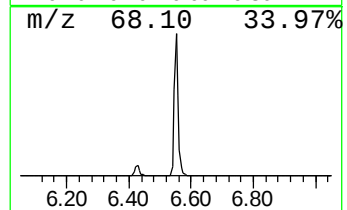
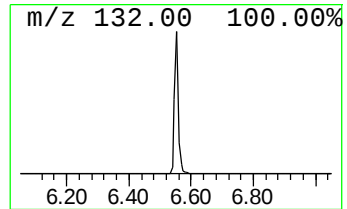
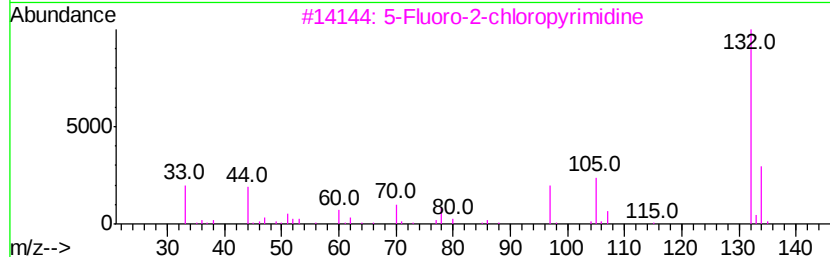
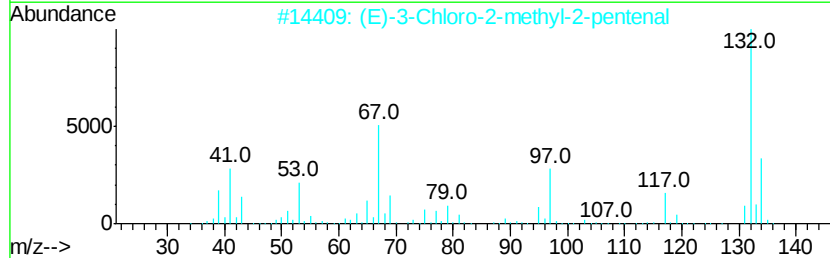
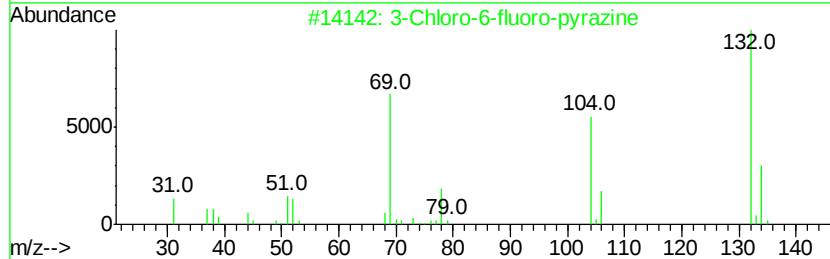
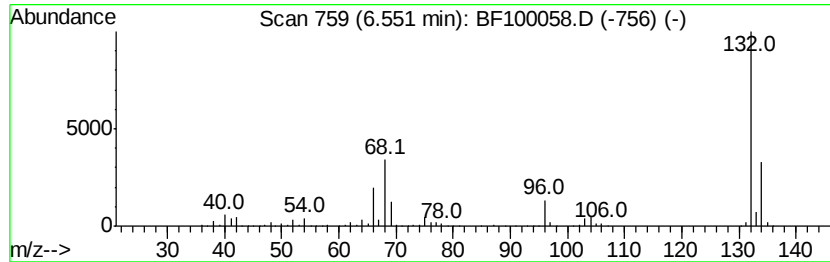
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.55 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.55	44.61 ng	1048540	1,4-Dichlorobenzene-d4	6.78

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	17
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
4		Xenon	132	Xe	007440-63-3	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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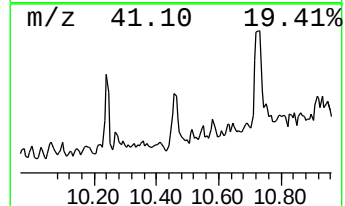
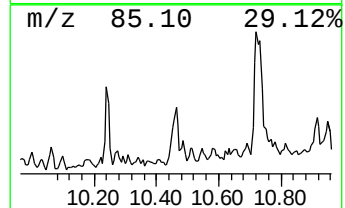
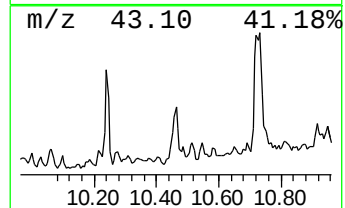
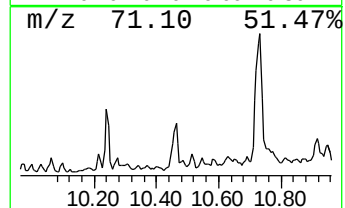
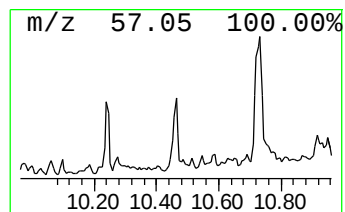
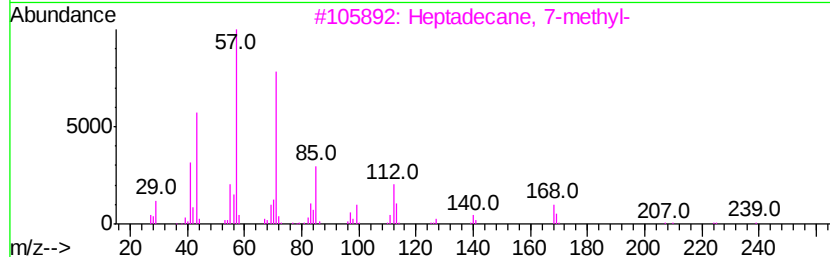
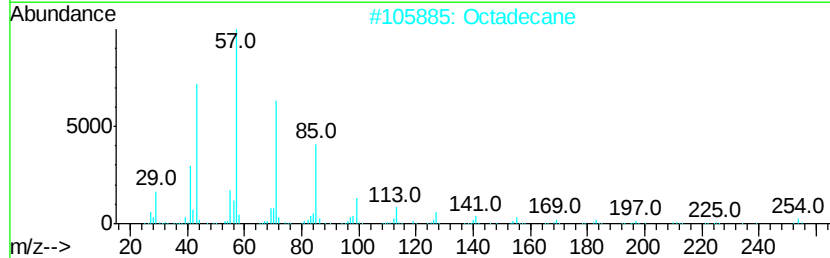
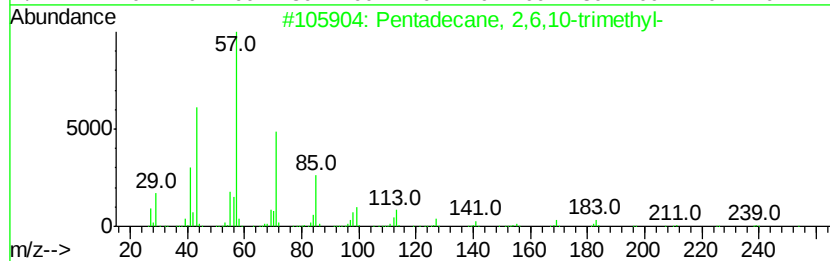
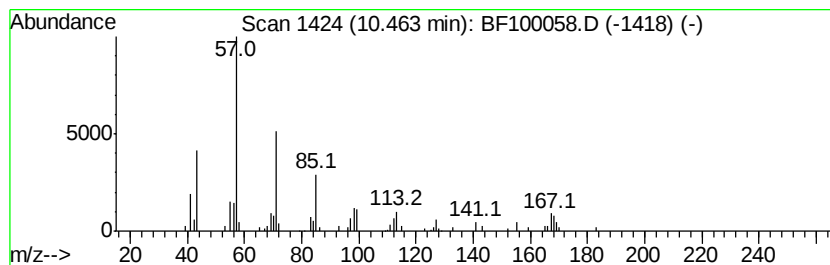
Instrument :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Pentadecane, 2,6,10-trimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.46	2.50 ng	79955	Acenaphthene-d10	9.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	80
2	Octadecane	254	C18H38	000593-45-3	72
3	Heptadecane, 7-methyl-	254	C18H38	020959-33-5	64
4	Eicosane	282	C20H42	000112-95-8	64
5	Hexacosane	366	C26H54	000630-01-3	62



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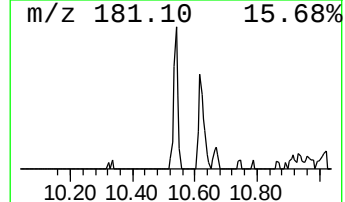
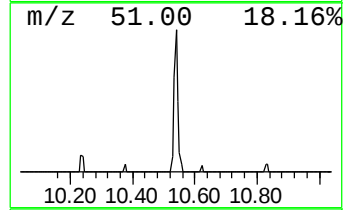
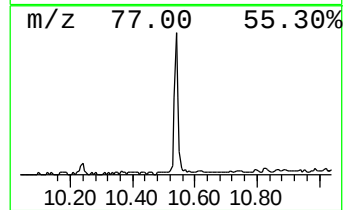
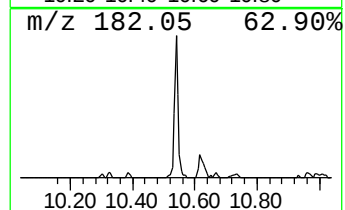
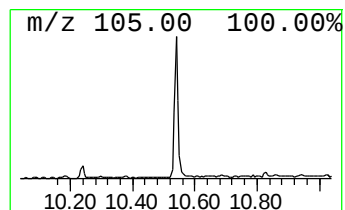
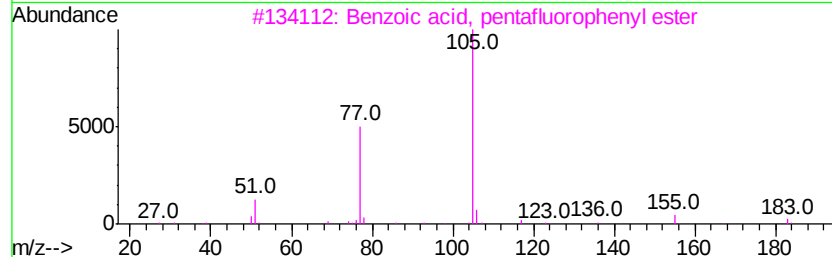
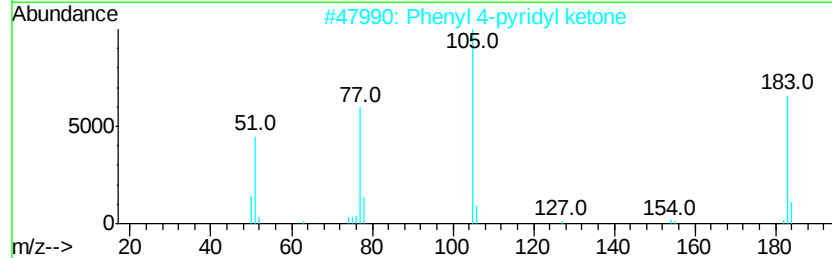
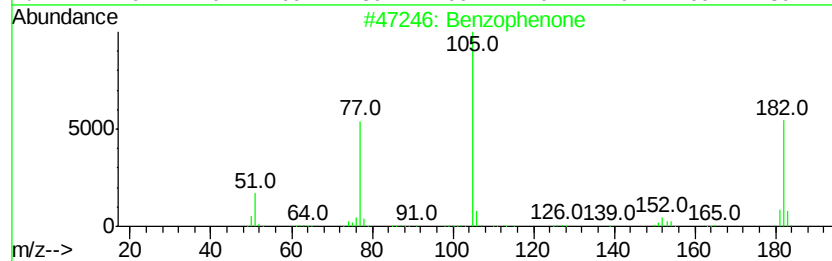
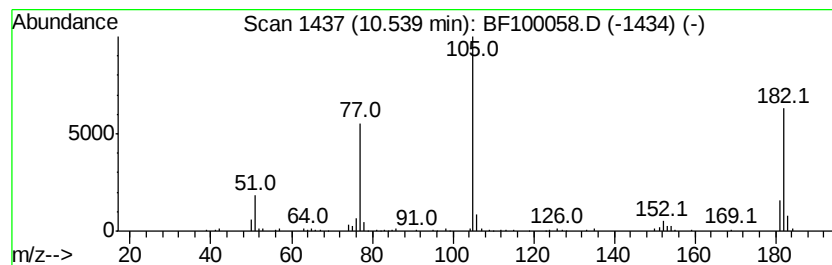
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzophenone Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.54	3.18 ng	101889	Acenaphthene-d10	9.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzophenone	182	C13H10O	000119-61-9	96
2		Phenyl 4-pyridyl ketone	183	C12H9NO	014548-46-0	47
3		Benzoic acid, pentafluorophenyl ...	288	C13H5F5O2	1000360-67-9	47
4		Phenylglyoxal	134	C8H6O2	001074-12-0	43
5		Benzoyl isothiocyanate	163	C8H5NOS	000532-55-8	43



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Data File : BF100058.D
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Operator : SJ/JU
Sample : I6236-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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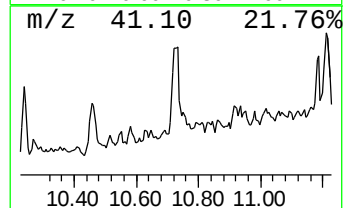
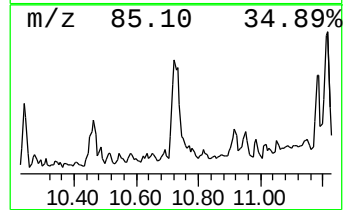
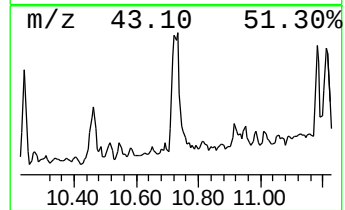
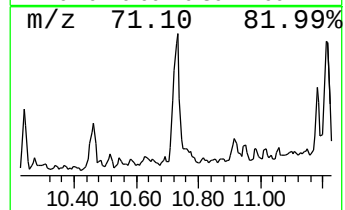
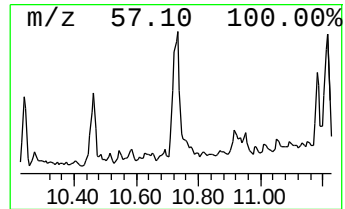
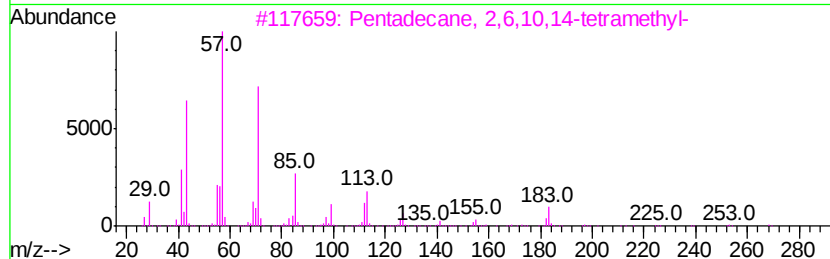
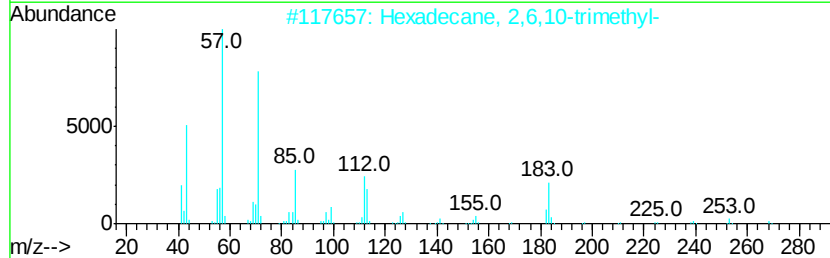
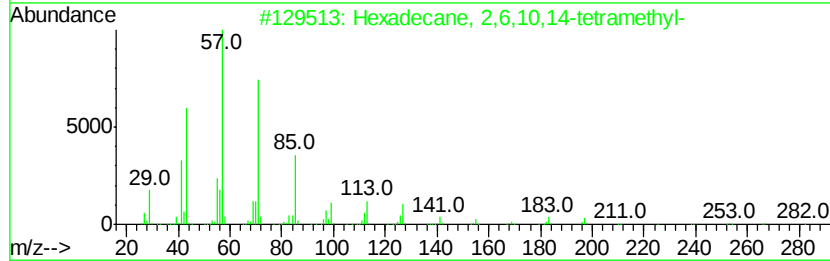
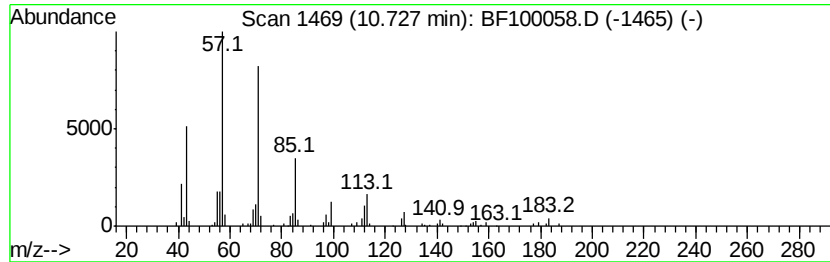
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Hexadecane, 2,6,10,14-tetra... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.73	7.69 ng	215891	Phenanthrene-d10	11.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	91
2		Hexadecane, 2,6,10-trimethyl-	268	C19H40	055000-52-7	86
3		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	80
4		Sulfurous acid, 2-ethylhexyl tri...	376	C21H44O3S	1000309-19-6	80
5		Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	80



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110117\
Data File : BF100058.D
Acq On : 1 Nov 2017 21:37
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ALS Vial : 16 Sample Multiplier: 1

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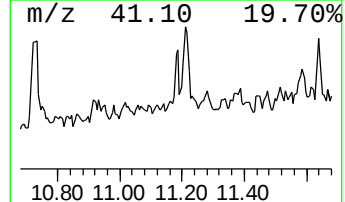
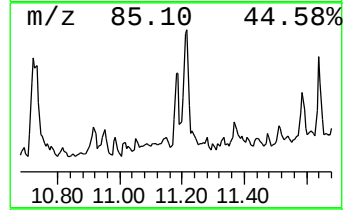
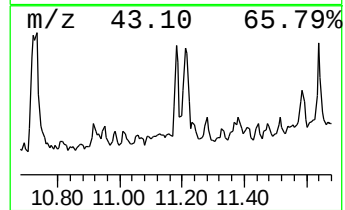
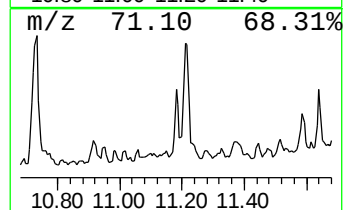
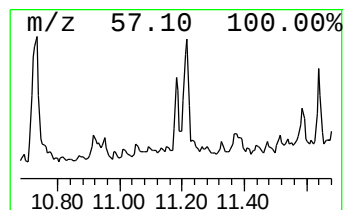
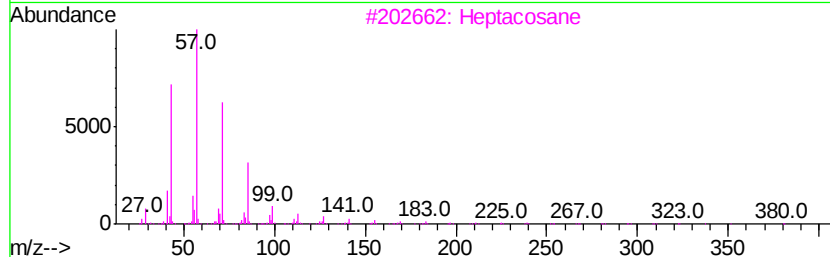
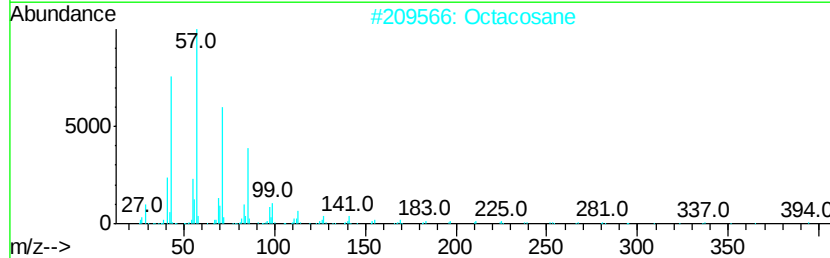
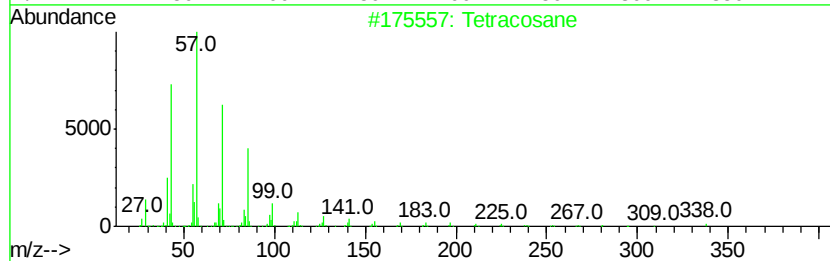
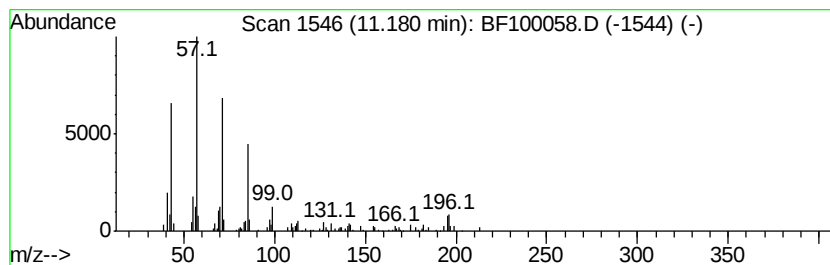
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tetracosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.18	2.55 ng	71694	Phenanthrene-d10	11.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	87
2		Octacosane	394	C28H58	000630-02-4	86
3		Heptacosane	380	C27H56	000593-49-7	86
4		Pentadecane	212	C15H32	000629-62-9	86
5		Heptadecane	240	C17H36	000629-78-7	80



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF110117\
Data File : BF100058.D
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Instrument :
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ClientSampleId :
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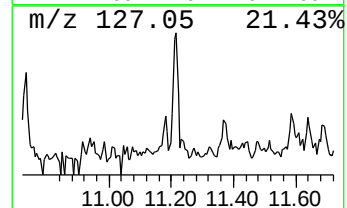
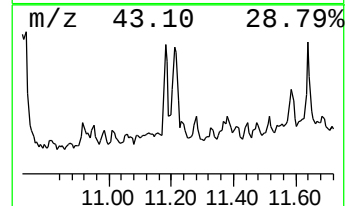
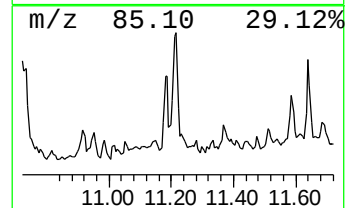
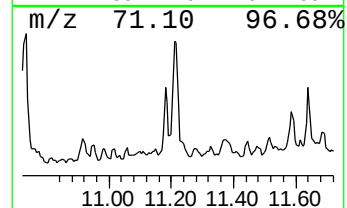
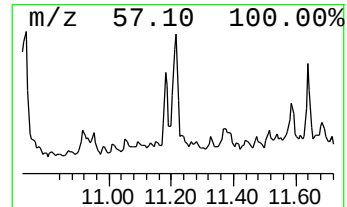
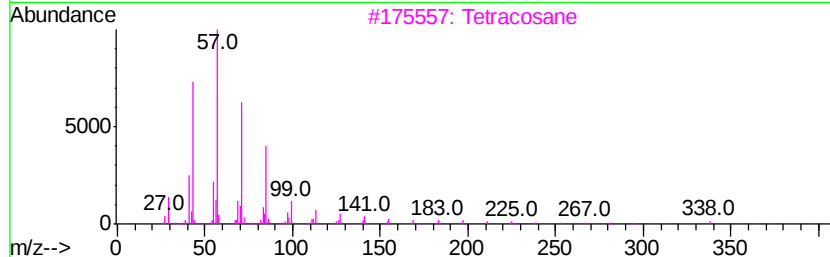
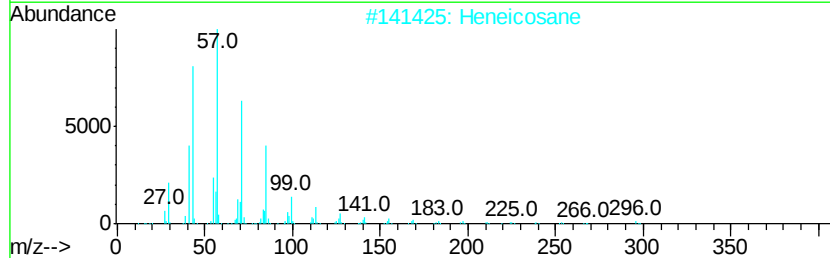
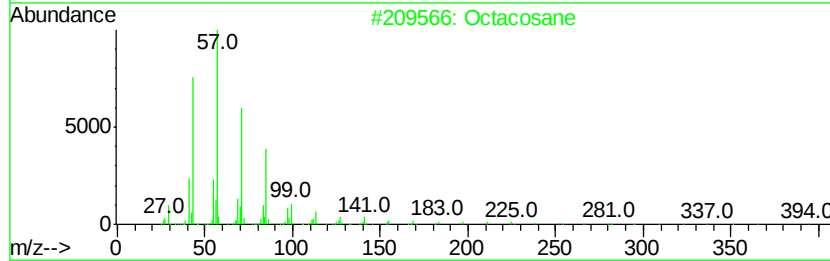
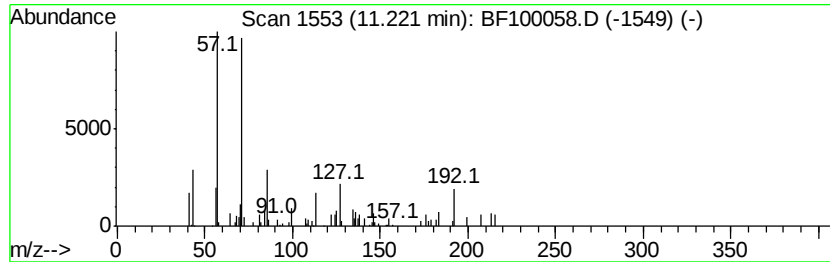
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Octacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.22	4.44 ng	124726	Phenanthrene-d10	11.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	86
2		Heneicosane	296	C21H44	000629-94-7	78
3		Tetracosane	338	C24H50	000646-31-1	78
4		Undecane, 3,8-dimethyl-	184	C13H28	017301-30-3	78
5		Decane, 3-methyl-	156	C11H24	013151-34-3	72



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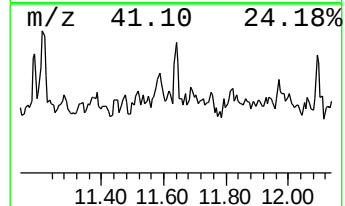
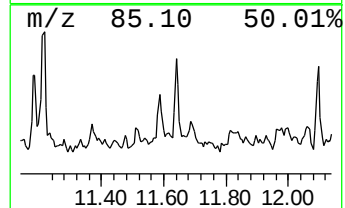
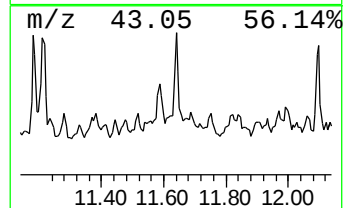
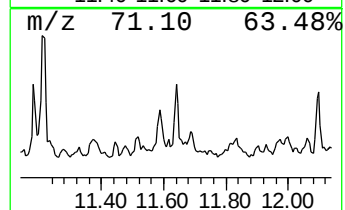
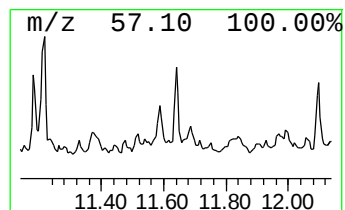
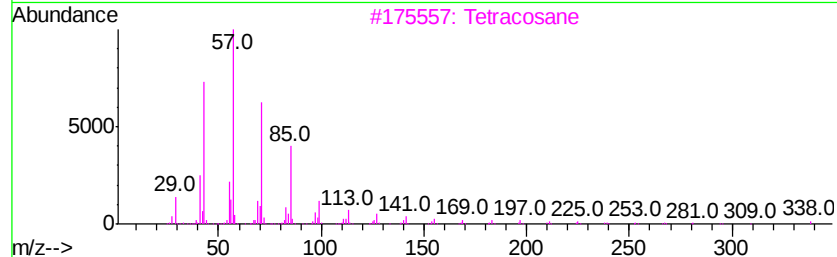
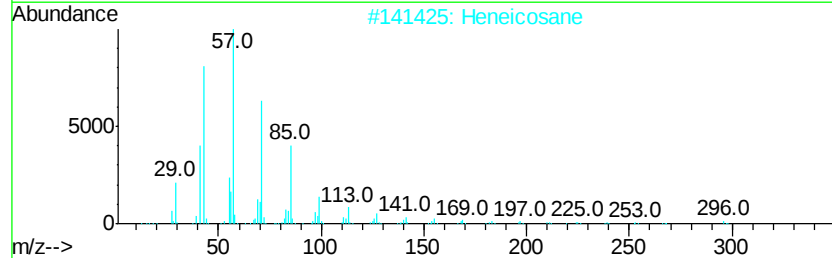
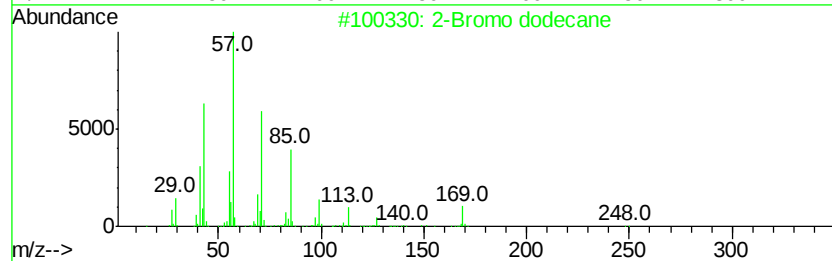
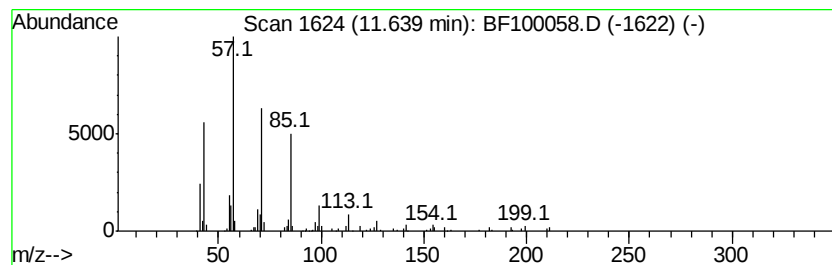
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 2-Bromo dodecane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.64	2.34 ng	65713	Phenanthrene-d10	11.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Bromo dodecane	248	C12H25Br	013187-99-0	90
2		Heneicosane	296	C21H44	000629-94-7	74
3		Tetracosane	338	C24H50	000646-31-1	72
4		Heptacosane	380	C27H56	000593-49-7	72
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	72



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 Data File : BF100058.D
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Instrument :
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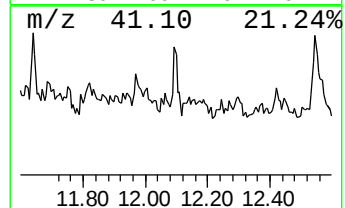
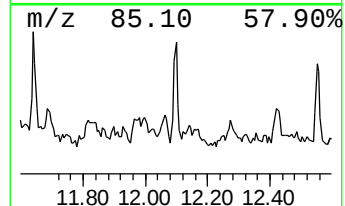
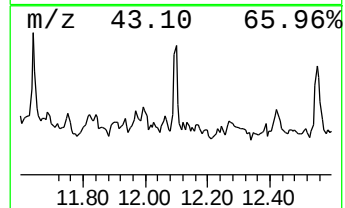
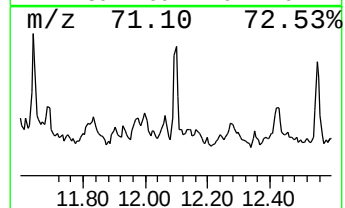
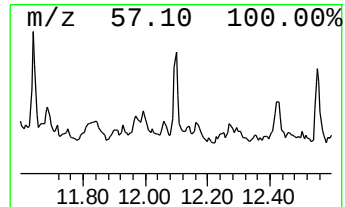
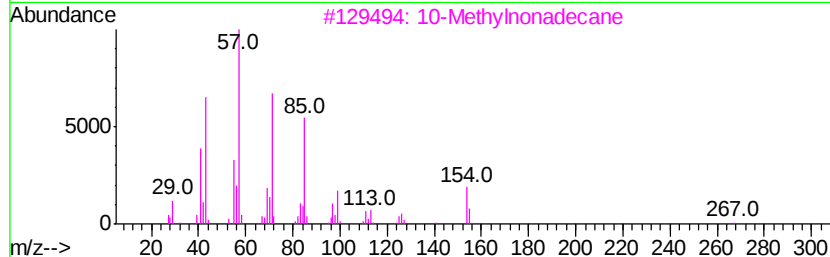
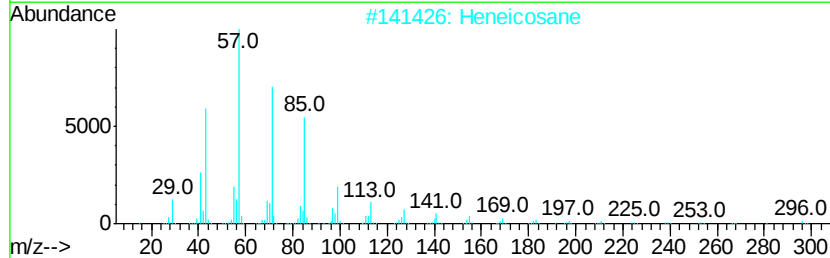
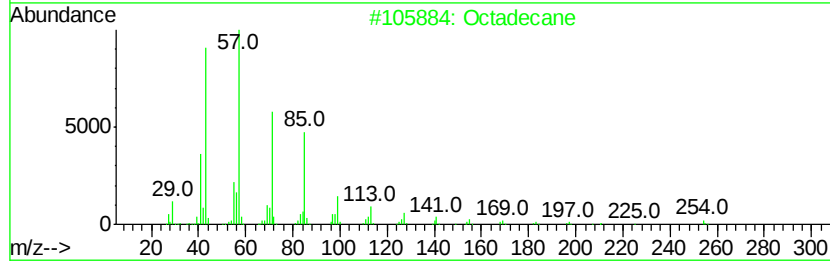
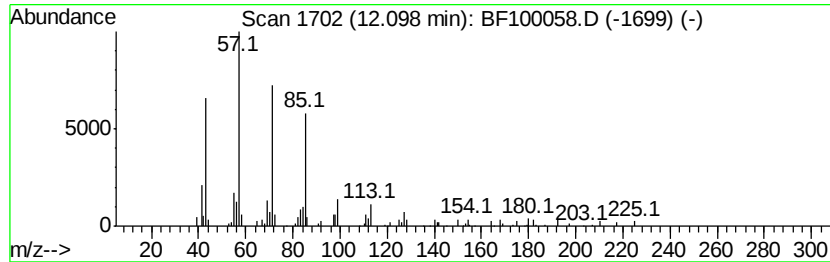
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 11 Octadecane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.10	2.36 ng	66374	Phenanthrene-d10	11.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	86
2		Heneicosane	296	C21H44	000629-94-7	86
3		10-Methylnonadecane	282	C20H42	056862-62-5	78
4		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	72
5		Undecane, 2,3-dimethyl-	184	C13H28	017312-77-5	64



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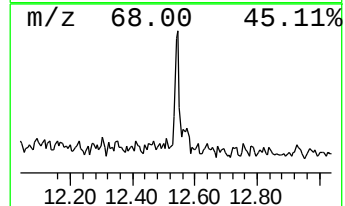
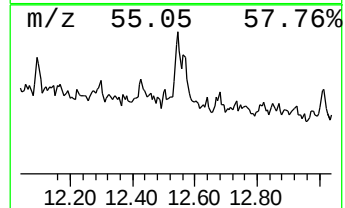
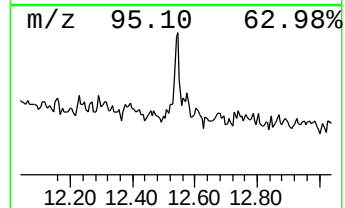
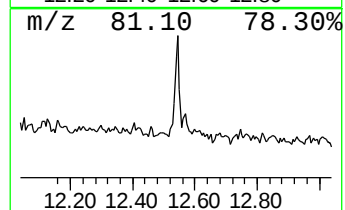
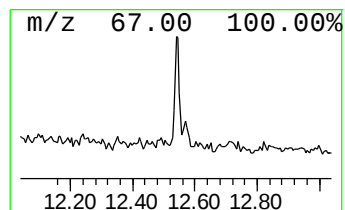
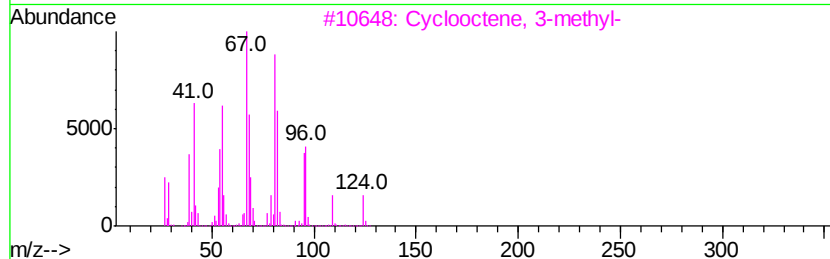
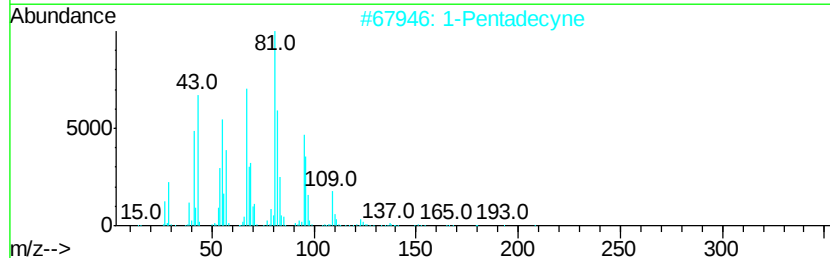
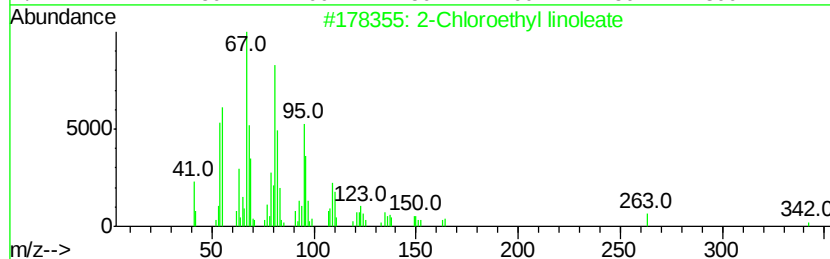
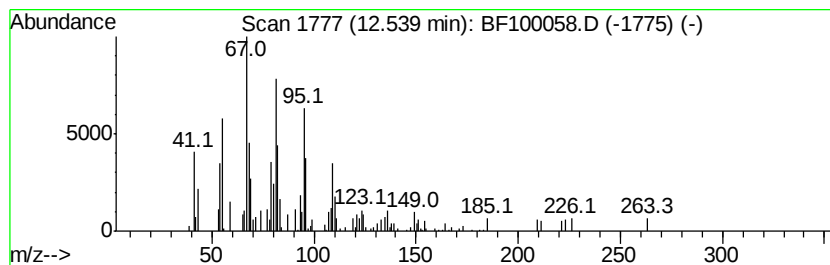
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 2-Chloroethyl linoleate Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	3.90 ng	109580	Phenanthrene-d10	11.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Chloroethyl linoleate	342	C20H35ClO2	025525-76-2	90
2		1-Pentadecyne	208	C15H28	000765-13-9	90
3		Cyclooctene, 3-methyl-	124	C9H16	013152-05-1	70
4		9,12-Octadecadienoic acid, methyl...	294	C19H34O2	002566-97-4	64
5		11,14-Eicosadienoic acid, methyl...	322	C21H38O2	002463-02-7	64



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Sample : I6236-01
Misc :
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Instrument :
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ClientSampleId :
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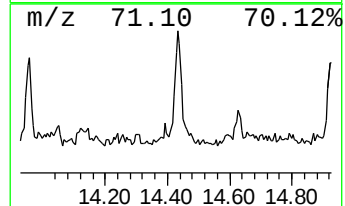
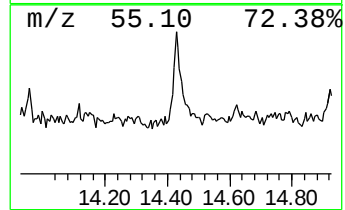
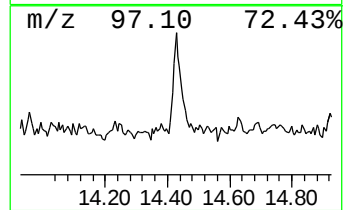
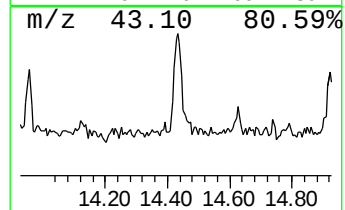
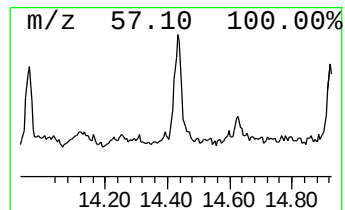
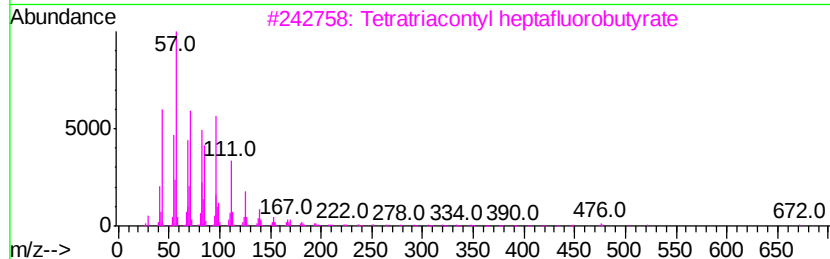
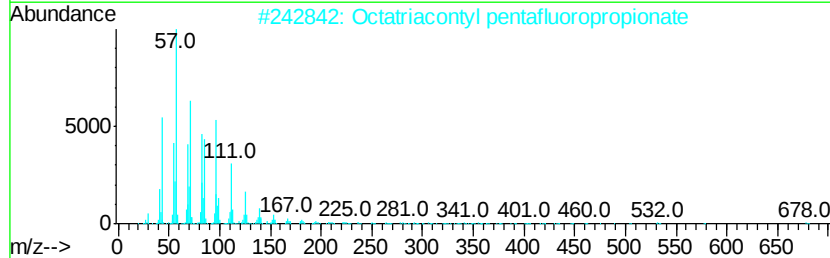
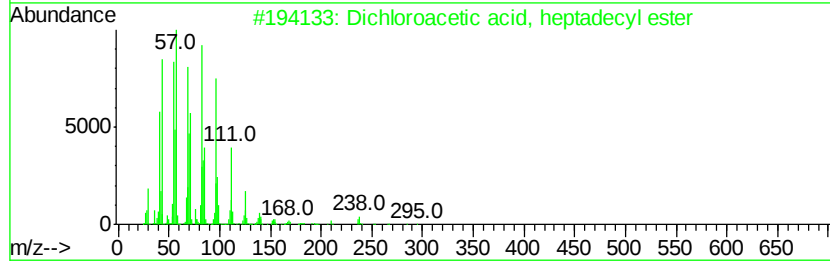
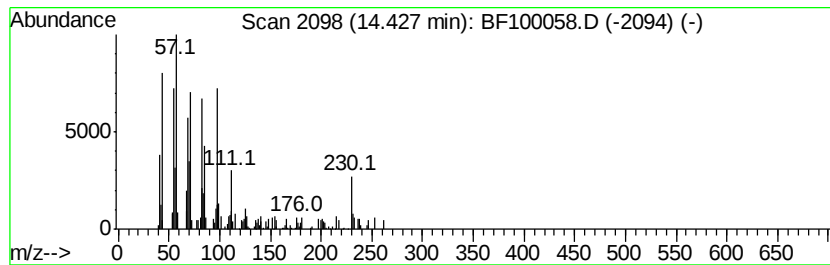
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Dichloroacetic acid, heptad... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.43	4.50 ng	150767	Chrysene-d12	14.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	53
2			Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	53
3			Tetratriacontyl heptafluorobutyrate	691	C38H69F7O2	1000351-84-1	53
4			Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	52
5			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	50



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110117\
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Sample : I6236-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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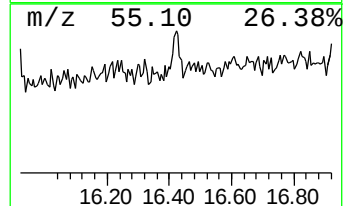
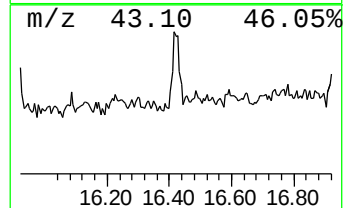
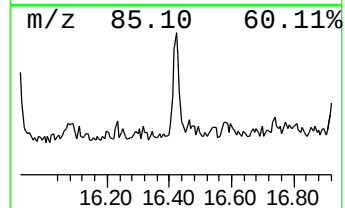
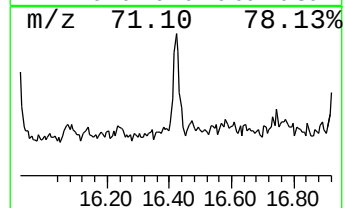
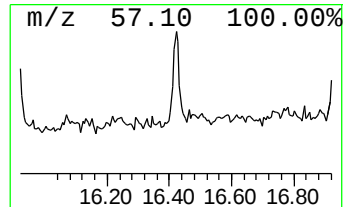
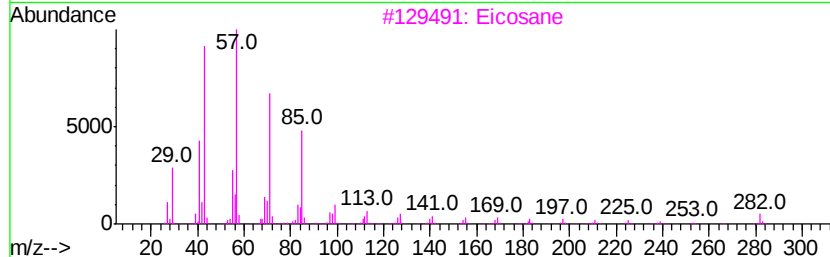
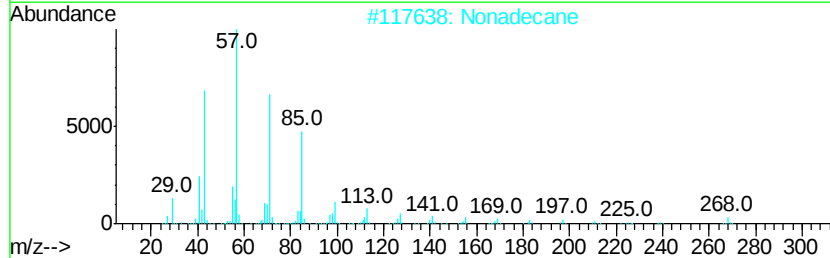
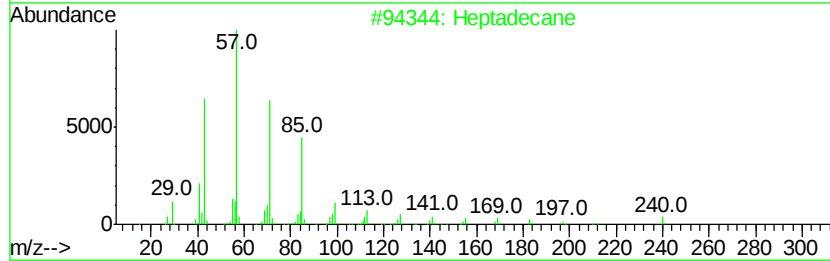
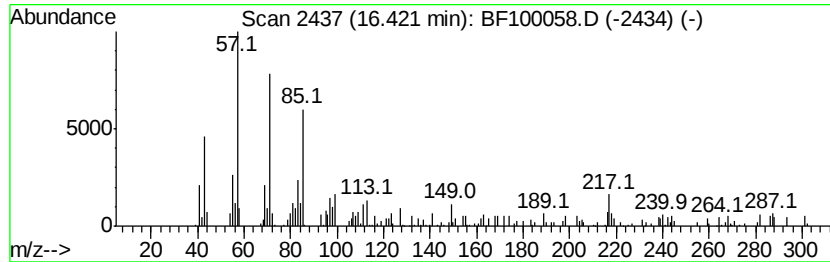
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heptadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.42	3.25 ng	64047	Perylene-d12	16.81

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	96
2	Nonadecane		268	C19H40	000629-92-5	93
3	Eicosane		282	C20H42	000112-95-8	89
4	Tetradecane		198	C14H30	000629-59-4	76
5	Heptacosane		380	C27H56	000593-49-7	68



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF110117\
Data File : BF100058.D
Acq On : 1 Nov 2017 21:37
Operator : SJ/JU
Sample : I6236-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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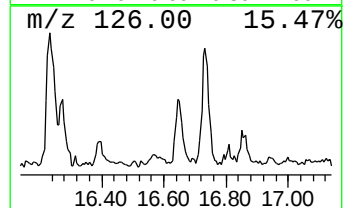
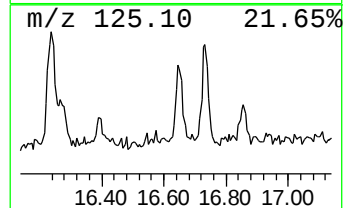
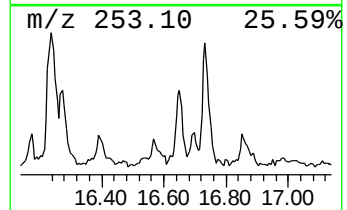
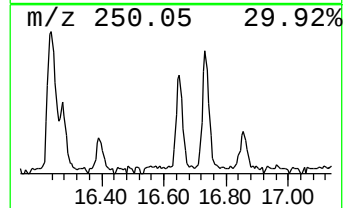
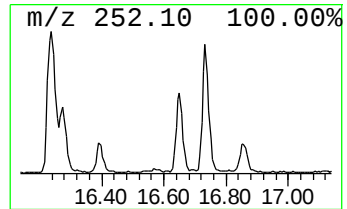
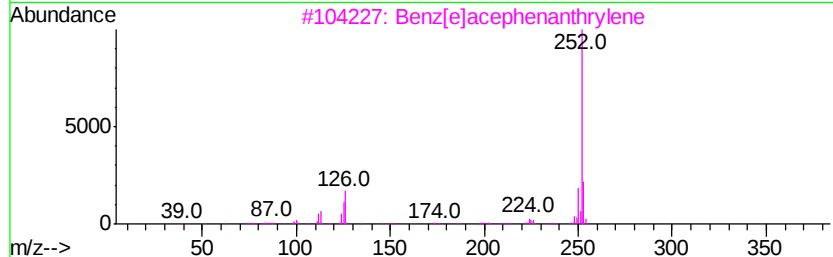
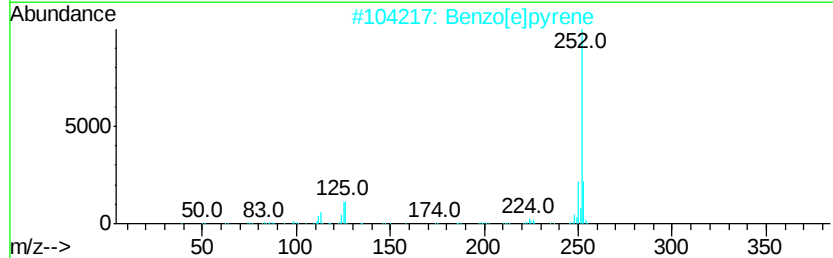
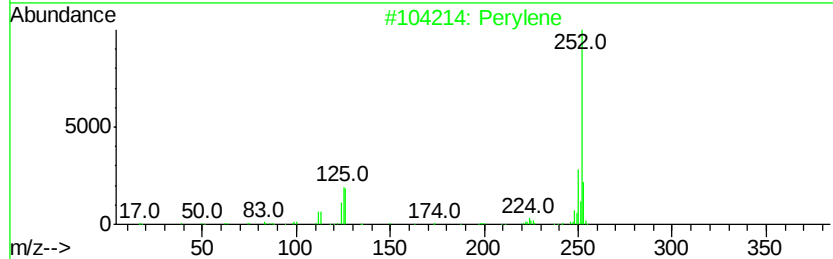
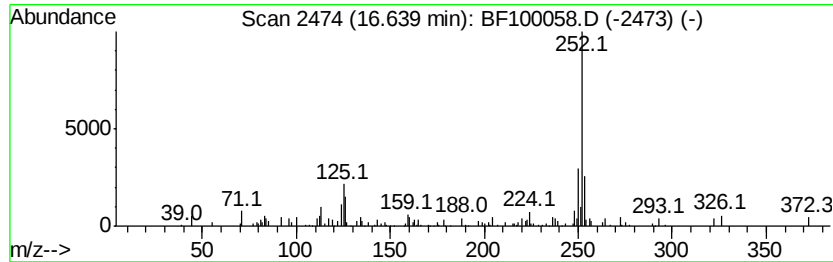
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Perylene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.64	3.93 ng	77498	Perylene-d12	16.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Perylene	252	C20H12	000198-55-0	96
2		Benzo[el]pyrene	252	C20H12	000192-97-2	94
3		Benzo[el]acephenanthrylene	252	C20H12	000205-99-2	90
4		Benzo[i]lfluoranthene	252	C20H12	000205-82-3	89
5		Benzo[k]fluoranthene	252	C20H12	000207-08-9	89



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110117\
Data File : BF100058.D
Acq On : 1 Nov 2017 21:37
Operator : SJ/JU
Sample : I6236-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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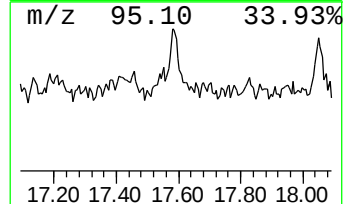
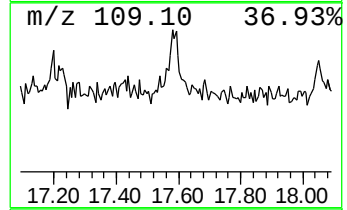
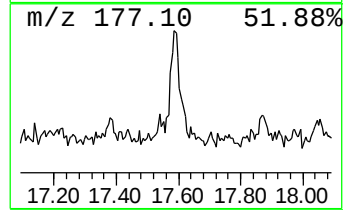
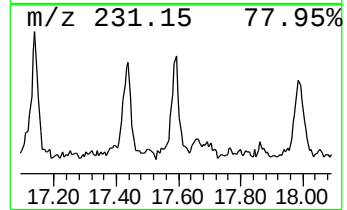
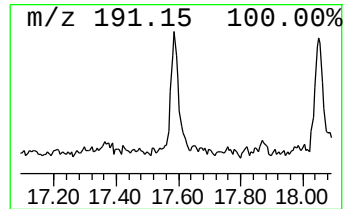
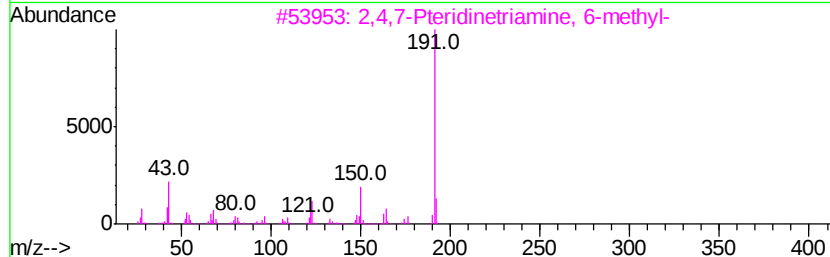
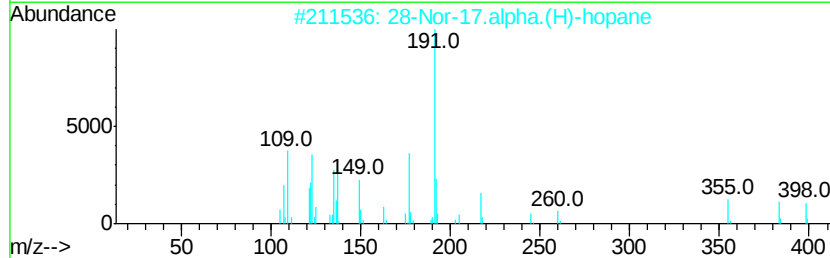
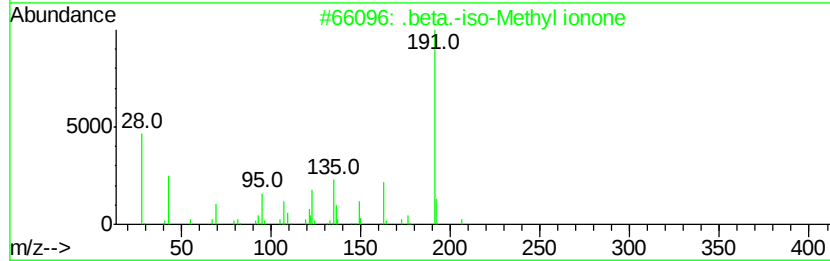
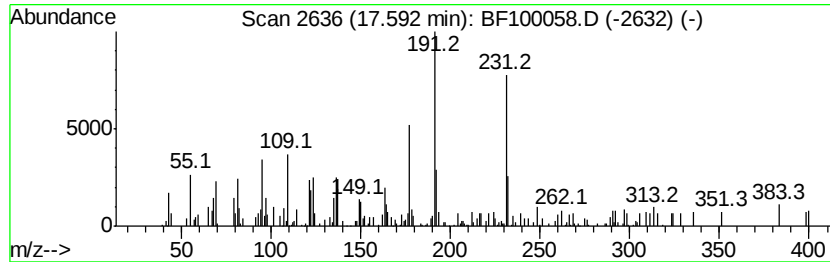
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 unknown17.59 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.59	6.34 ng	125002	Perylene-d12	16.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	46
2			28-Nor-17.alpha.(H)-hopane	398	C29H50	053584-60-4	38
3			2,4,7-Pteridinetriamine, 6-methyl-	191	C7H9N7	017539-50-3	38
4			Anthracene, 1-(3-butenyl)-	232	C18H16	1000156-51-8	27
5			Anthracene, 9-(3-butenyl)-	232	C18H16	097451-55-3	27



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF110117\
 Data File : BF100058.D
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 Operator : SJ/JU
 Sample : I6236-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one, 4...	4.35	16.0	ng	375785	1	6.78	470141	20.0
2-Pentanone, 4-hy...	4.99	47.3	ng	1112940	1	6.78	470141	20.0
unknown6.55	6.55	44.6	ng	1048540	1	6.78	470141	20.0
Pentadecane, 2,6,...	10.46	2.5	ng	79955	3	9.82	640220	20.0
Benzophenone	10.54	3.2	ng	101889	3	9.82	640220	20.0
Hexadecane, 2,6,1...	10.73	7.7	ng	215891	4	11.35	561396	20.0
Tetracosane	11.18	2.5	ng	71694	4	11.35	561396	20.0
Octacosane	11.22	4.4	ng	124726	4	11.35	561396	20.0
2-Bromo dodecane	11.64	2.3	ng	65713	4	11.35	561396	20.0
Octadecane	12.10	2.4	ng	66374	4	11.35	561396	20.0
2-Chloroethyl lin...	12.54	3.9	ng	109580	4	11.35	561396	20.0
Dichloroacetic ac...	14.43	4.5	ng	150767	5	14.63	670365	20.0
Heptadecane	16.42	3.3	ng	64047	6	16.81	394333	20.0
Perylene	16.64	3.9	ng	77498	6	16.81	394333	20.0
unknown17.59	17.59	6.3	ng	125002	6	16.81	394333	20.0