

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081717\
 Data File : BF097810.D
 Acq On : 18 Aug 2017 2:34
 Operator : SJ/JU
 Sample : I4774-02
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HY-SB441

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-BF081517.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	3.787	284	289	295	rVB	88461	119911	3.33%	0.369%
2	4.975	487	491	500	rVB	367764	488717	13.59%	1.505%
3	5.351	550	555	556	rBV	61314	53870	1.50%	0.166%
4	5.387	556	561	564	rVB	3336253	3488028	97.01%	10.740%
5	5.604	593	598	606	rVB2	56563	90506	2.52%	0.279%
6	6.369	724	728	730	rBV2	30947	42209	1.17%	0.130%
7	6.410	730	735	738	rVV	3281661	3595699	100.00%	11.072%
8	6.545	753	758	761	rBV	3397624	3356806	93.36%	10.336%
9	6.610	765	769	773	rVB3	106580	118400	3.29%	0.365%
10	6.775	793	797	801	rBV	997293	820659	22.82%	2.527%
11	6.934	820	824	827	rBV	2999439	2566876	71.39%	7.904%
12	7.034	838	841	843	rBV	54641	48181	1.34%	0.148%
13	7.339	885	893	896	rBV	2273589	2129334	59.22%	6.557%
14	7.381	897	900	904	rVV3	48071	55560	1.55%	0.171%
15	7.422	904	907	911	rVB4	45318	50422	1.40%	0.155%
16	7.810	969	973	976	rBV3	32484	40792	1.13%	0.126%
17	8.057	1010	1015	1017	rBV	1103617	981791	27.30%	3.023%
18	8.081	1017	1019	1022	rVB	532738	399651	11.11%	1.231%
19	8.657	1113	1117	1119	rBV2	54874	49349	1.37%	0.152%
20	8.704	1119	1125	1130	rVV9	18293	46267	1.29%	0.142%
21	8.769	1133	1136	1140	rVB	381332	308144	8.57%	0.949%
22	8.869	1149	1153	1156	rBV	171952	134486	3.74%	0.414%
23	9.086	1187	1190	1194	rVB	45400	41535	1.16%	0.128%
24	9.139	1194	1199	1202	rBV	3281853	2995714	83.31%	9.224%
25	9.233	1208	1215	1218	rBV2	116910	188079	5.23%	0.579%
26	9.328	1229	1231	1236	rVB	45547	52529	1.46%	0.162%
27	9.398	1239	1243	1247	rVV2	84183	131529	3.66%	0.405%
28	9.469	1252	1255	1258	rVV	96285	97235	2.70%	0.299%
29	9.498	1258	1260	1262	rVV	75347	56124	1.56%	0.173%
30	9.528	1262	1265	1270	rVV	541814	499972	13.90%	1.540%
31	9.586	1273	1275	1279	rVB	41106	45285	1.26%	0.139%
32	9.675	1286	1290	1294	rBV	150050	171001	4.76%	0.527%
33	9.751	1301	1303	1307	rBV	84908	68989	1.92%	0.212%
34	9.810	1307	1313	1317	rBV	826032	833873	23.19%	2.568%

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35	10.016	1345	1348	1351	rVB2	67373	64938	1.81%	0.200%
36	10.251	1386	1388	1395	rVB2	77192	92503	2.57%	0.285%
37	10.357	1403	1406	1410	rBV4	52489	84081	2.34%	0.259%
38	10.475	1423	1426	1429	rVB2	58023	57867	1.61%	0.178%
39	10.598	1443	1447	1450	rBV	1491381	1311069	36.46%	4.037%
40	10.722	1466	1468	1477	rVB2	126409	222643	6.19%	0.686%
41	11.298	1562	1566	1568	rBV	726735	686647	19.10%	2.114%
42	11.316	1568	1569	1575	rVV	161238	132579	3.69%	0.408%
43	11.369	1576	1578	1581	rVB	108194	82796	2.30%	0.255%
44	11.510	1600	1602	1604	rBV	55517	46871	1.30%	0.144%
45	11.598	1615	1617	1620	rVB2	59483	38260	1.06%	0.118%
46	11.827	1653	1656	1661	rVB4	104045	126215	3.51%	0.389%
47	12.345	1742	1744	1747	rVB	72691	61597	1.71%	0.190%
48	12.504	1769	1771	1774	rVB	136175	118404	3.29%	0.365%
49	12.733	1808	1810	1812	rBV	245927	177286	4.93%	0.546%
50	12.763	1812	1815	1819	rVB	198790	171169	4.76%	0.527%
51	12.886	1831	1836	1839	rBV	2655723	2380323	66.20%	7.330%
52	13.168	1881	1884	1886	rBV	74985	66540	1.85%	0.205%
53	13.257	1897	1899	1902	rVB	114380	93496	2.60%	0.288%
54	13.286	1902	1904	1908	rBV	76339	64194	1.79%	0.198%
55	13.786	1986	1989	1996	rVB	207555	234205	6.51%	0.721%
56	13.927	2009	2013	2016	rBV2	860150	945644	26.30%	2.912%
57	14.951	2184	2187	2195	rVB	102257	195231	5.43%	0.601%
58	15.239	2233	2236	2242	rVB	193557	215932	6.01%	0.665%
59	15.298	2243	2246	2252	rVB	155514	187111	5.20%	0.576%
60	15.362	2253	2257	2260	rBV	383921	450570	12.53%	1.387%

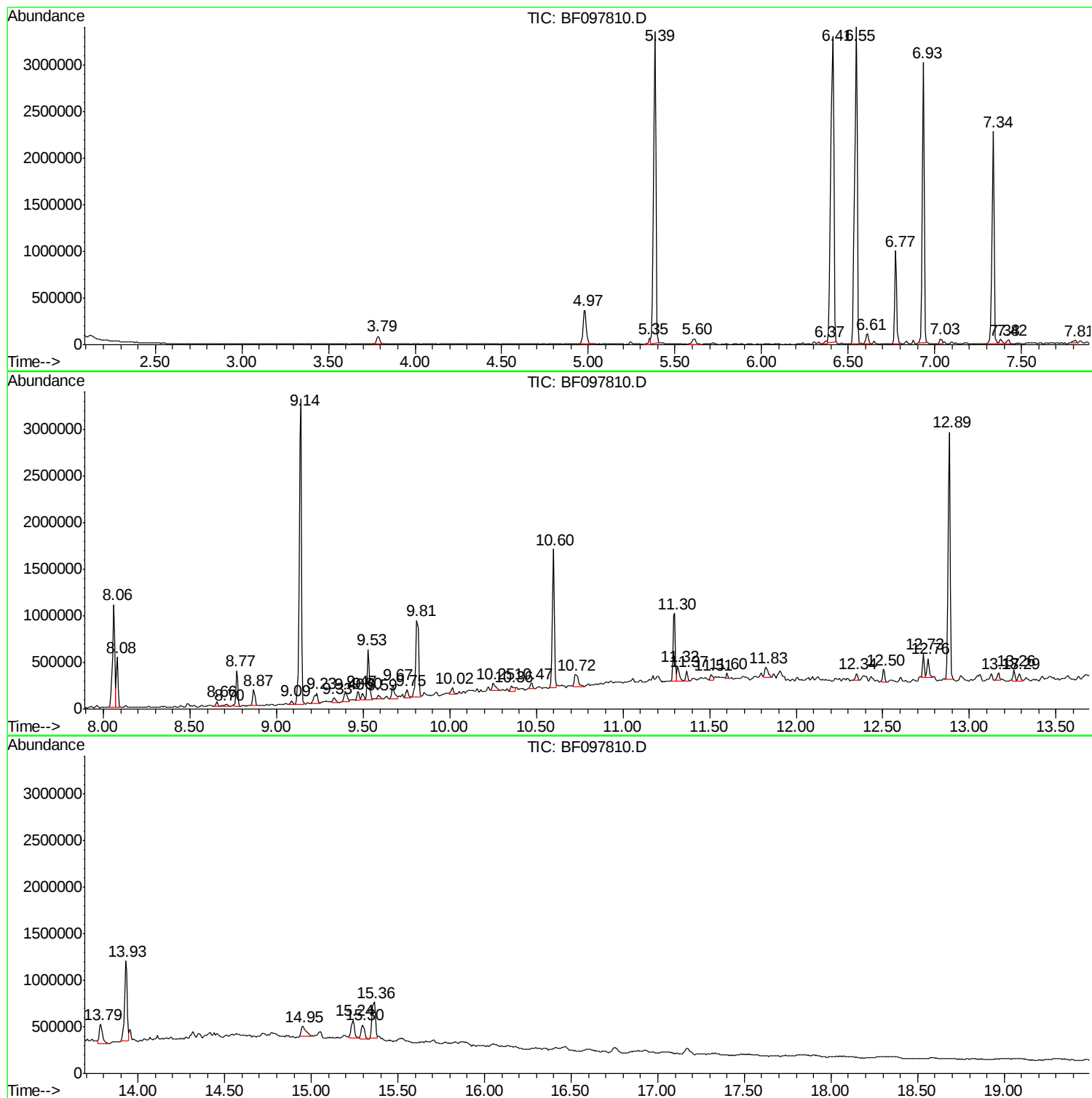
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TIC Library : C:\DATABASE\NIST11.L
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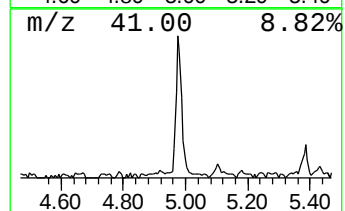
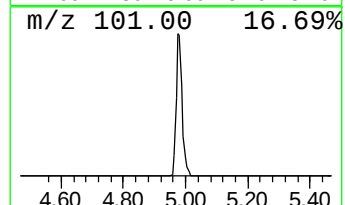
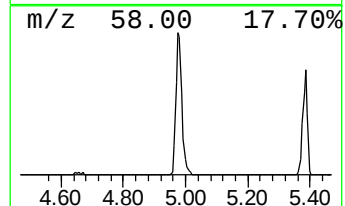
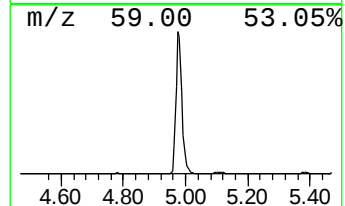
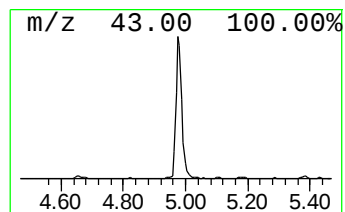
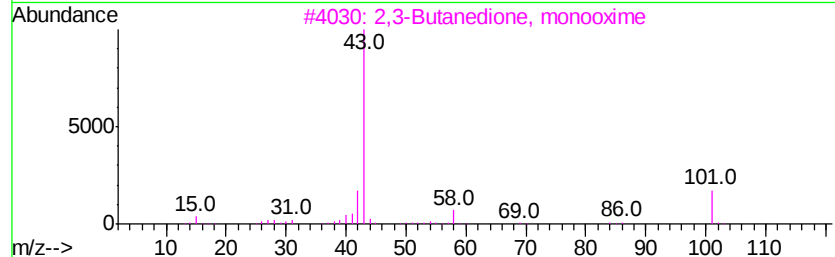
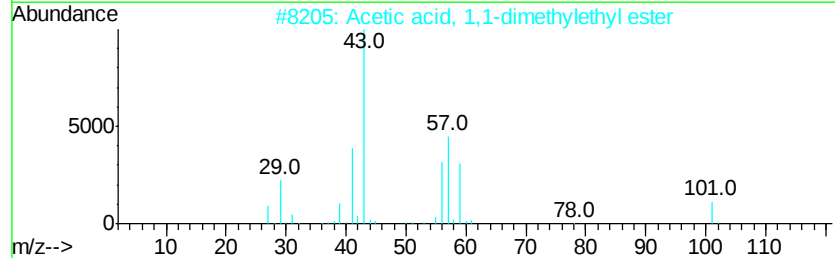
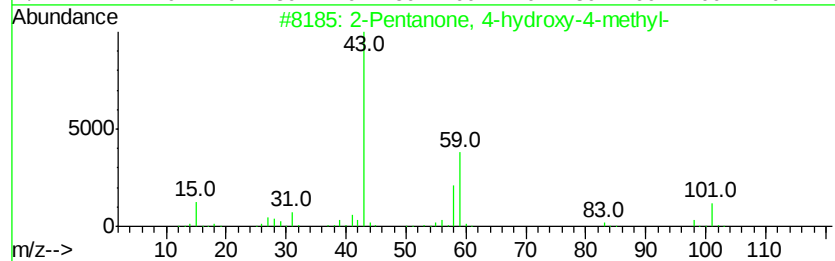
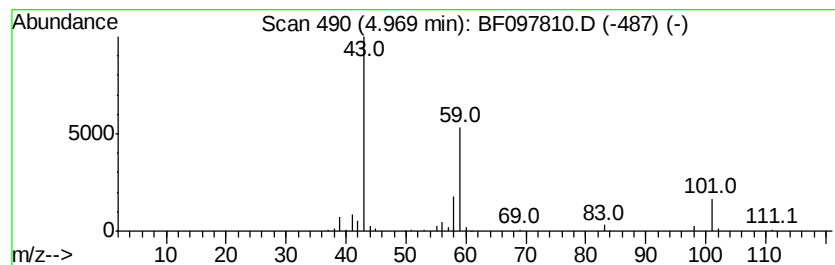
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.97	11.91 ng	488717	1,4-Dichlorobenzene-d4	6.77

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	28
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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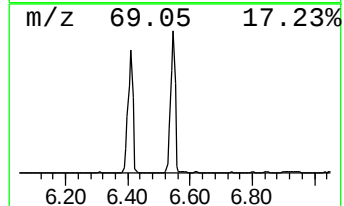
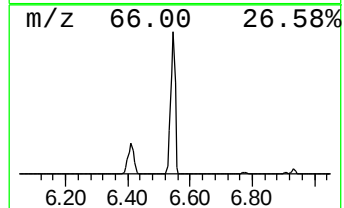
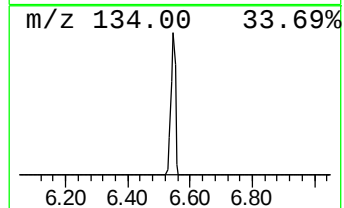
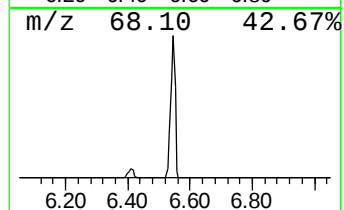
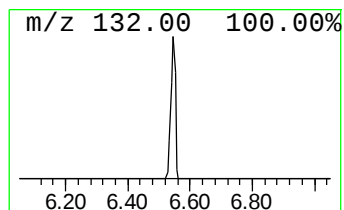
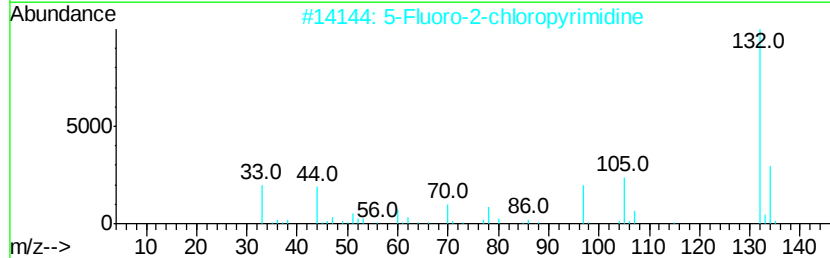
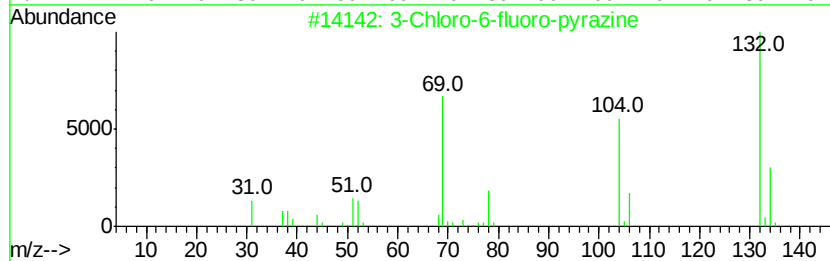
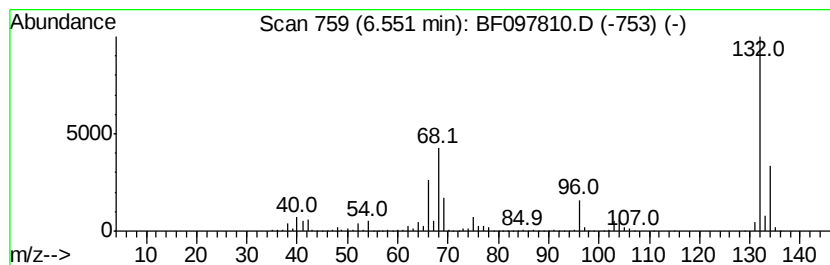
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 unknown6.55 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.55	81.81 ng	3356810	1,4-Dichlorobenzene-d4	6.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3			5-Aminoindole	132	C8H8N2	005192-03-0	10
4			1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5			Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	9



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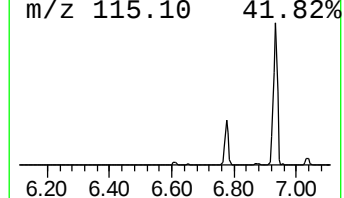
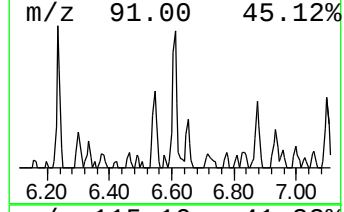
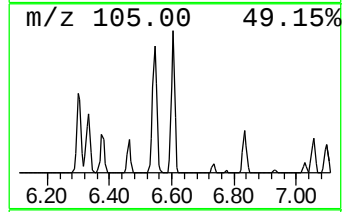
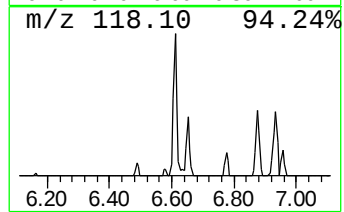
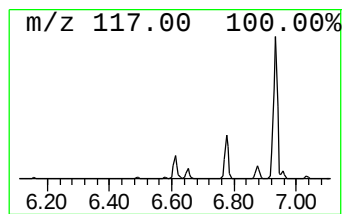
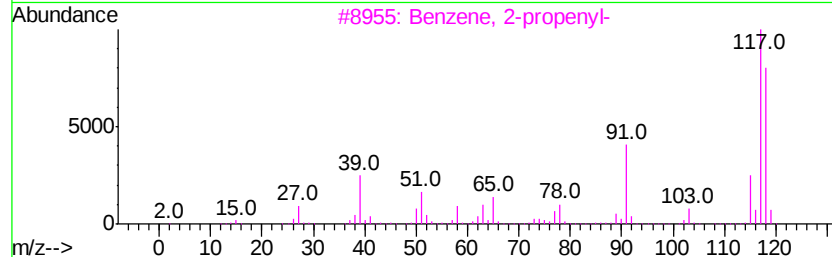
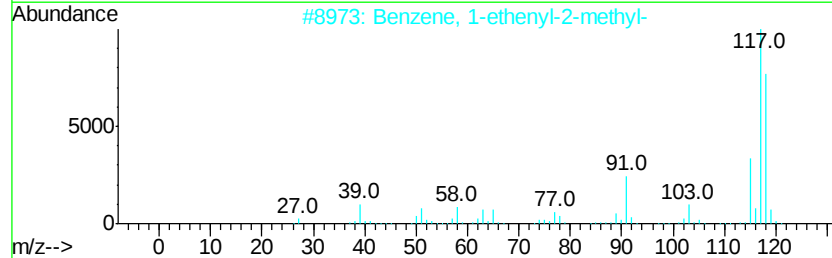
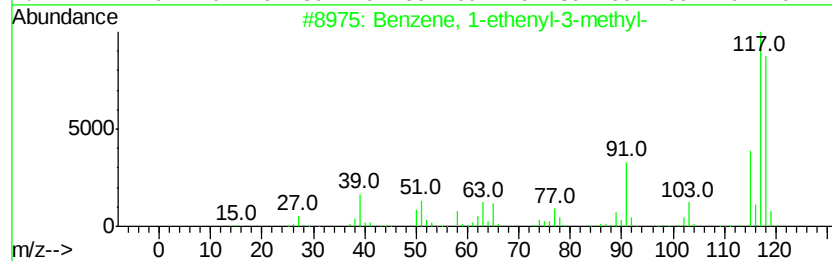
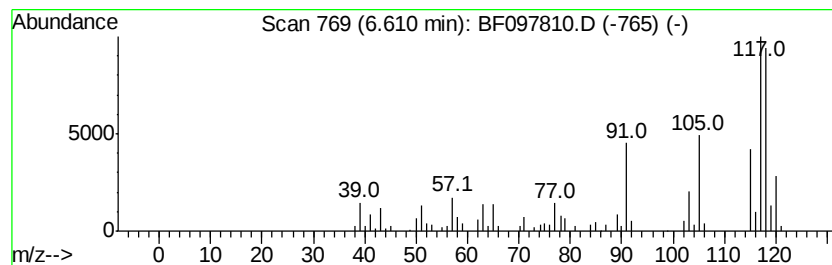
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-ethenyl-3-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	2.89 ng	118400	1,4-Dichlorobenzene-d4	6.77

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	93
2		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	92
3		Benzene, 2-propenyl-	118	C9H10	000300-57-2	91
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	022250-74-4	86
5		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	86



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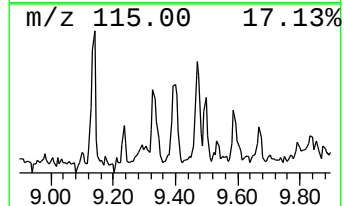
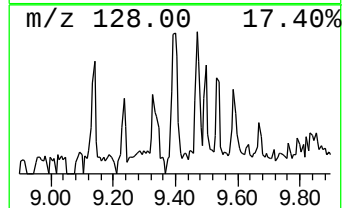
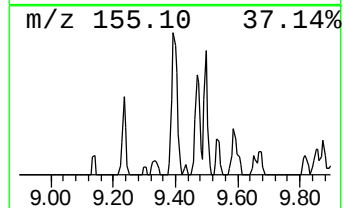
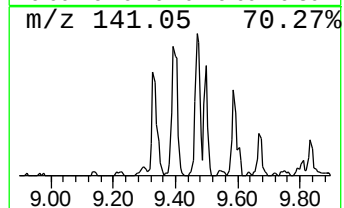
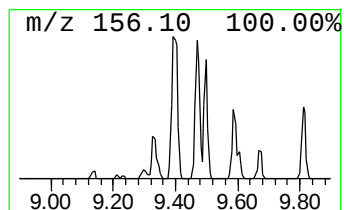
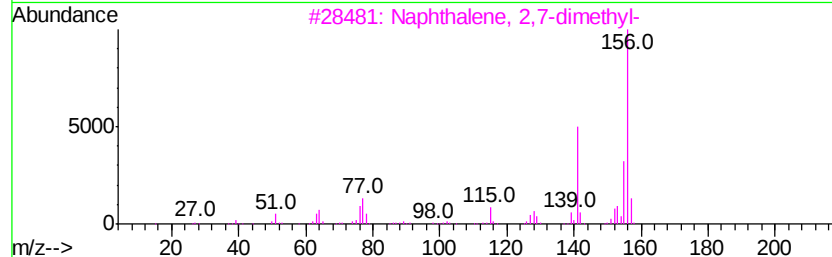
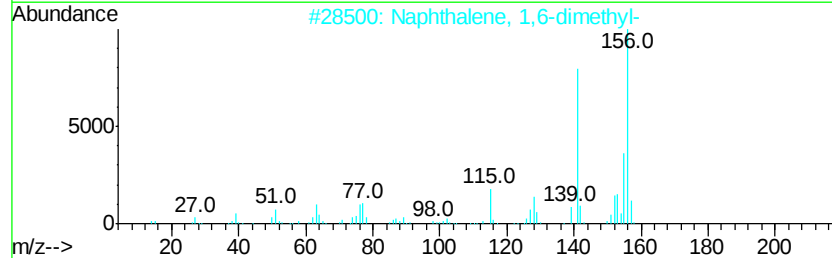
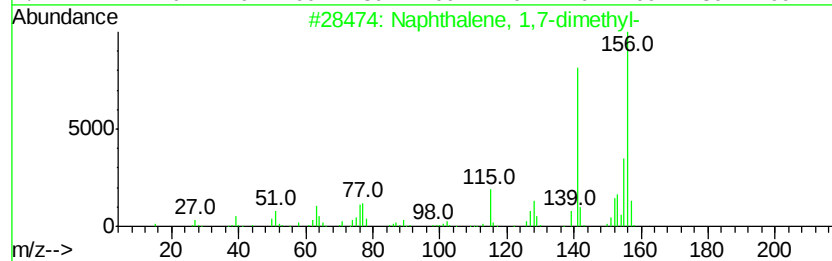
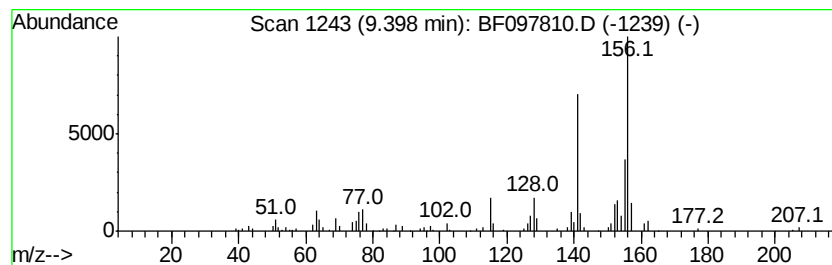
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Naphthalene, 1,7-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.40	3.15 ng	131529	Acenaphthene-d10	9.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
2		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
3		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
5		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97



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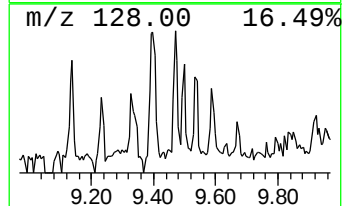
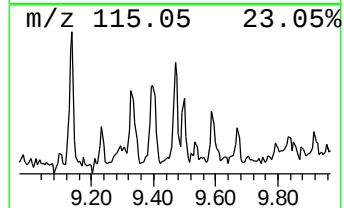
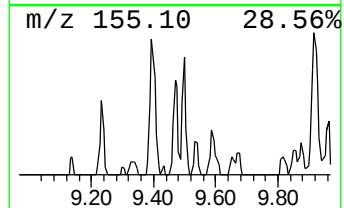
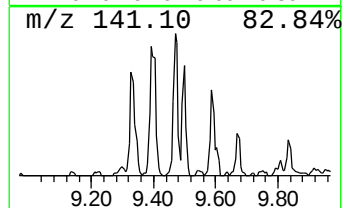
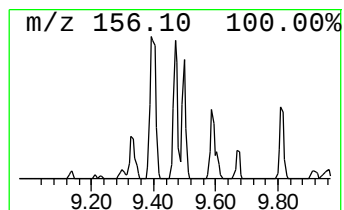
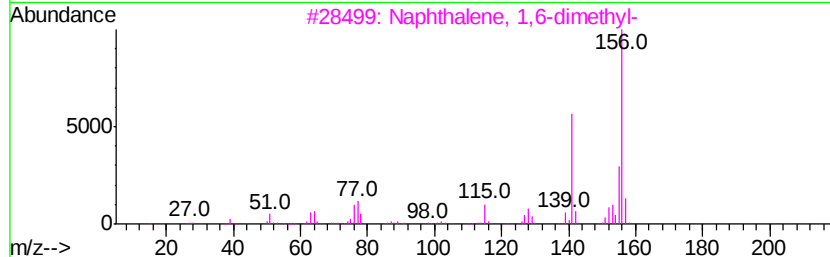
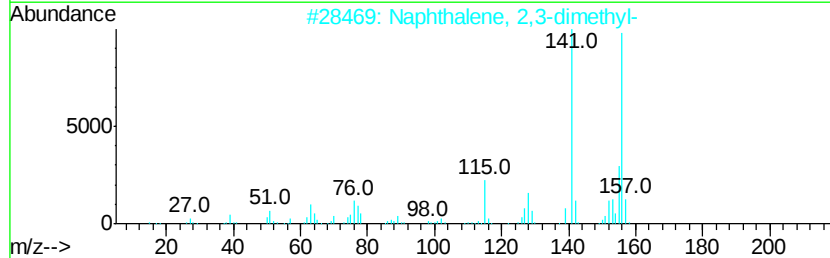
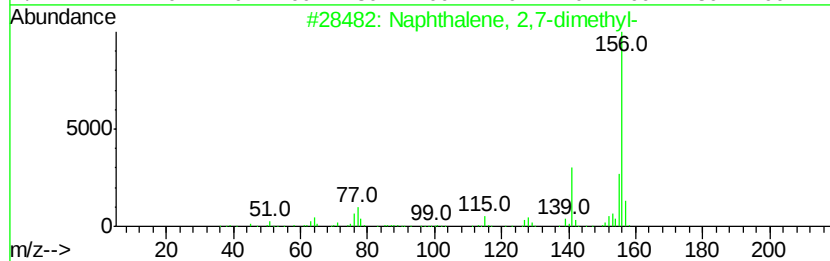
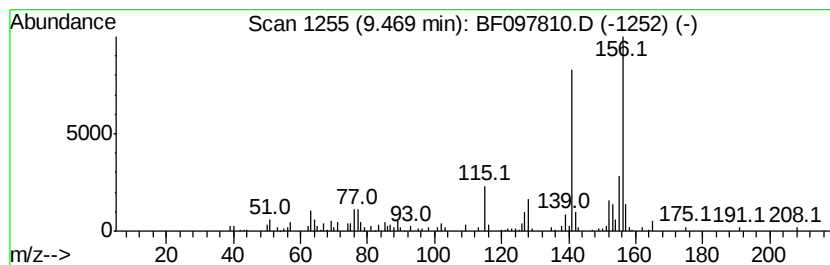
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Naphthalene, 2,7-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.47	2.33 ng	97235	Acenaphthene-d10	9.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
2		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
3		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	94
4		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	94
5		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081717\
Data File : BF097810.D
Acq On : 18 Aug 2017 2:34
Operator : SJ/JU
Sample : I4774-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

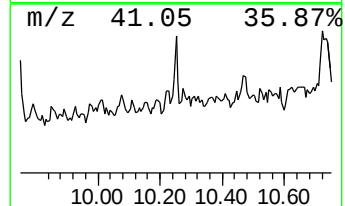
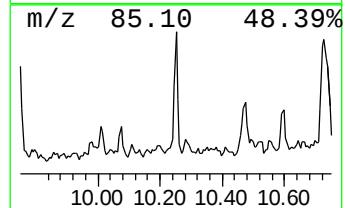
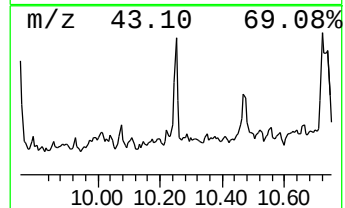
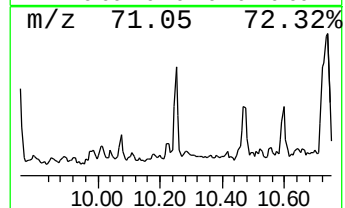
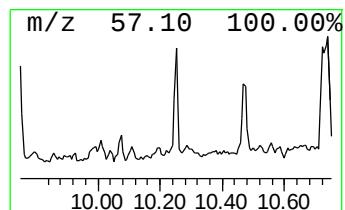
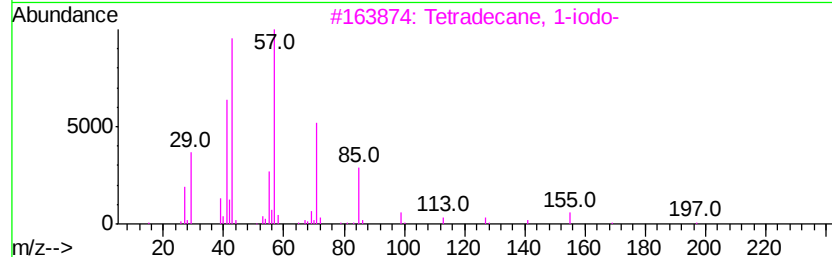
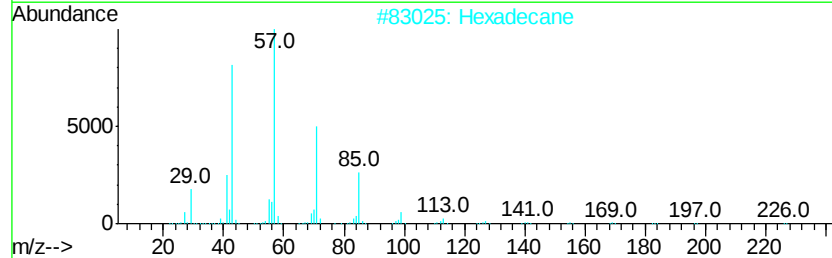
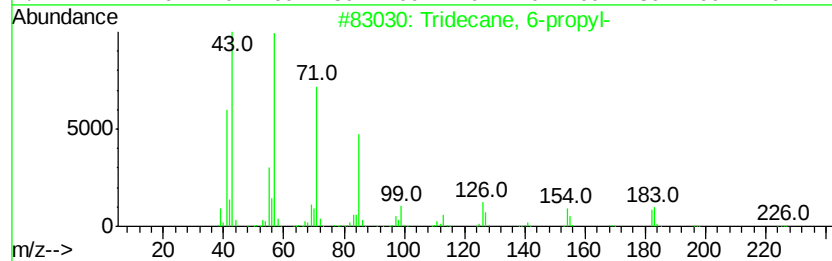
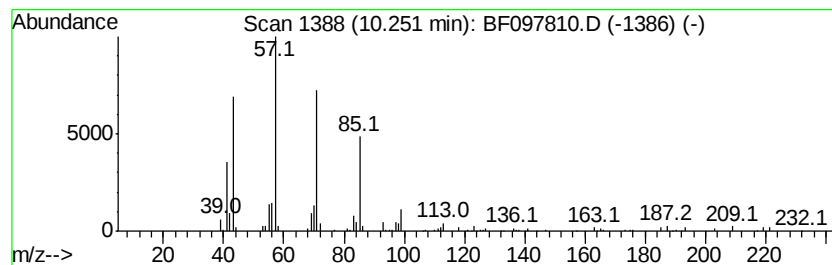
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tridecane, 6-propyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.25	2.22 ng	92503	Acenaphthene-d10	9.81
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tridecane, 6-propyl-	226 C16H34	055045-10-8 83
2		Hexadecane	226 C16H34	000544-76-3 80
3		Tetradecane, 1-iodo-	324 C14H29I	019218-94-1 72
4		Dodecane, 2-methyl-6-propyl-	226 C16H34	055045-08-4 72
5		Octadecane, 1-iodo-	380 C18H37I	000629-93-6 64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081717\
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Operator : SJ/JU
Sample : I4774-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

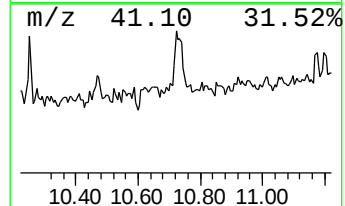
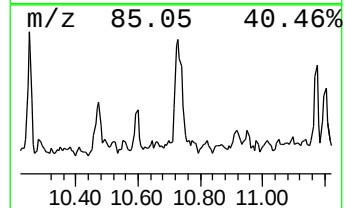
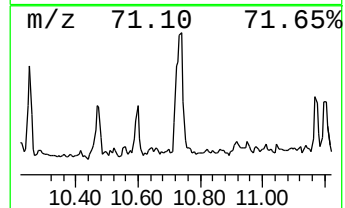
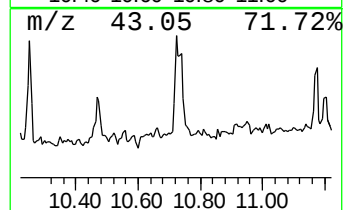
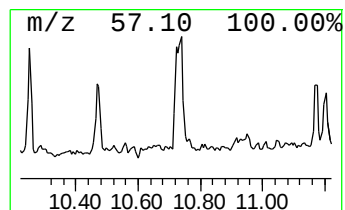
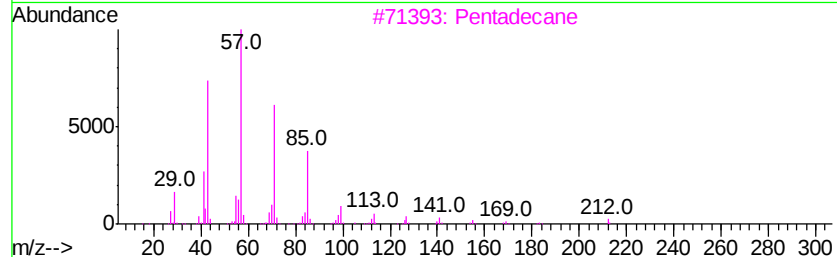
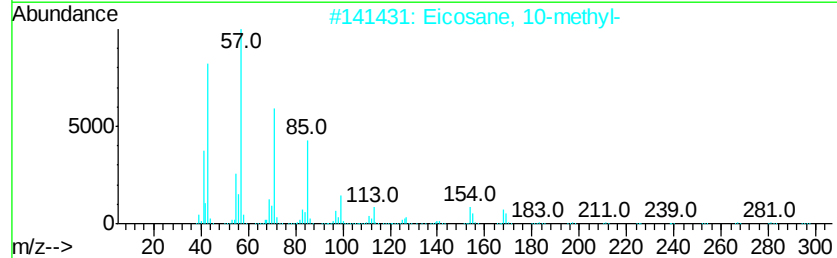
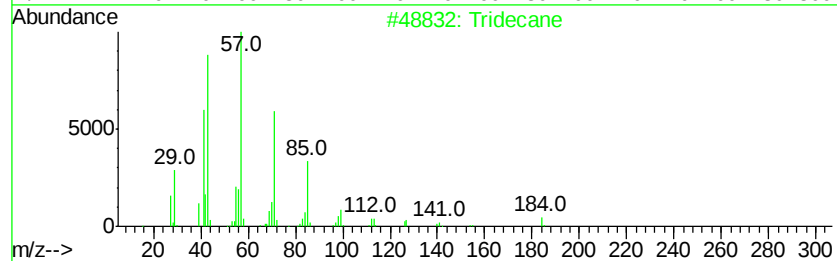
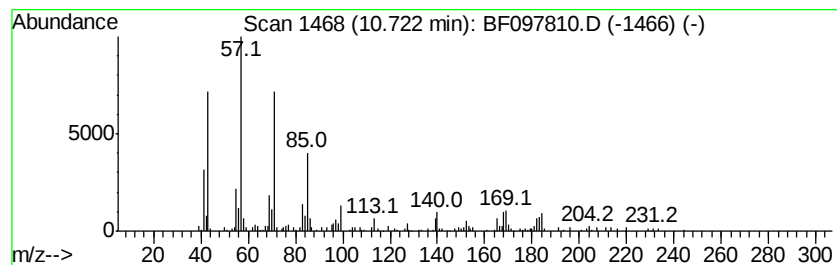
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tridecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.72	6.48 ng	222643	Phenanthrene-d10		11.30
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	74
2	Eicosane, 10-methyl-	296	C21H44	054833-23-7	72
3	Pentadecane	212	C15H32	000629-62-9	64
4	Eicosane	282	C20H42	000112-95-8	64
5	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081717\
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Sample : I4774-02
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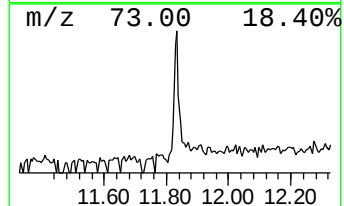
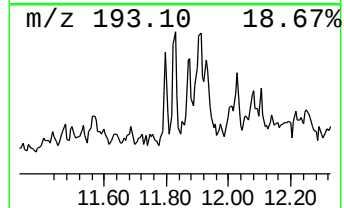
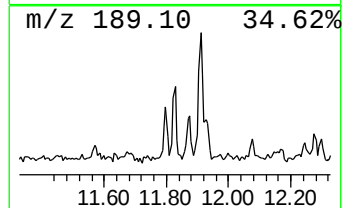
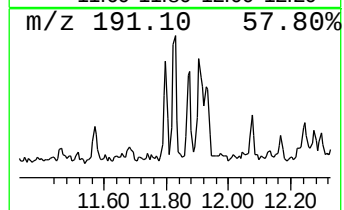
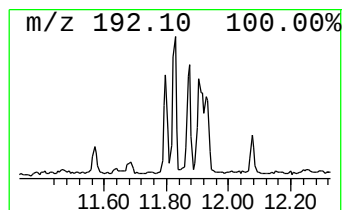
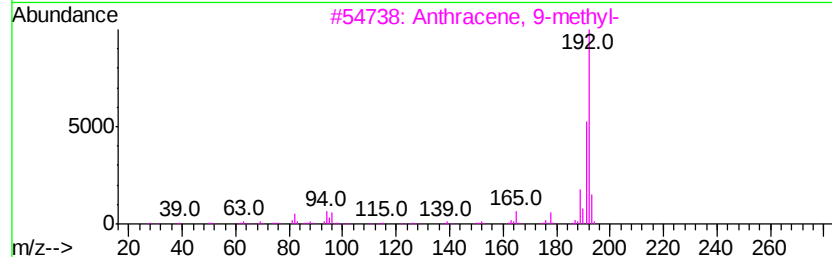
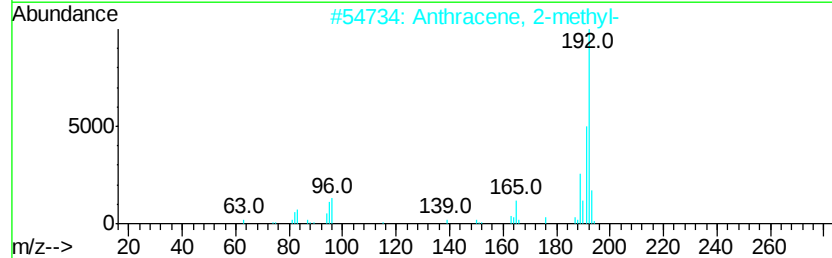
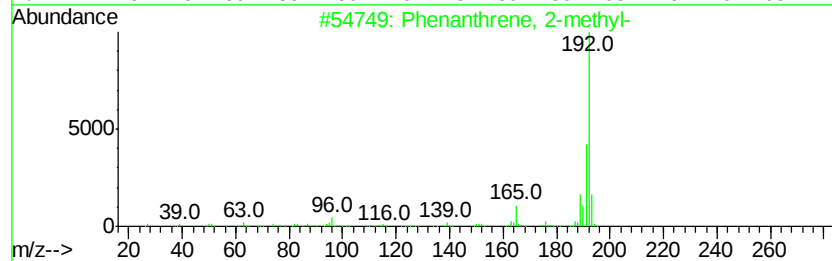
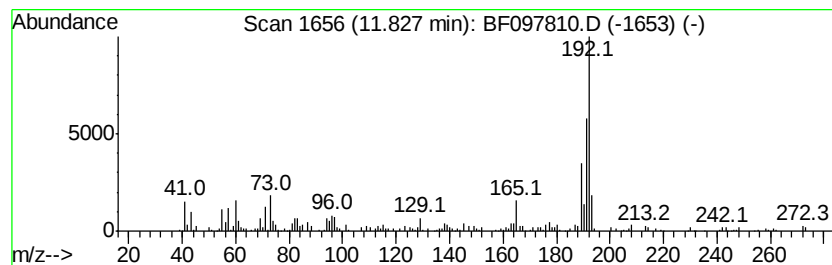
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Phenanthrene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.83	3.68 ng	126215	Phenanthrene-d10	11.30	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	95
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
5	1H-Cyclopropa[1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081717\
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Sample : I4774-02
Misc :
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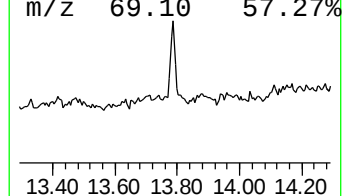
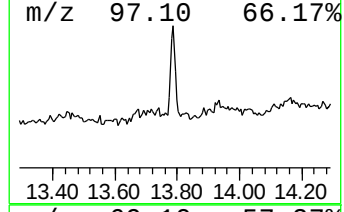
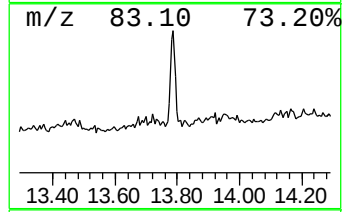
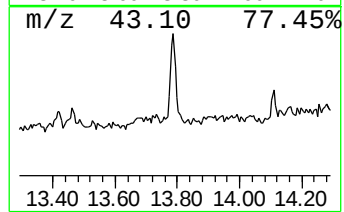
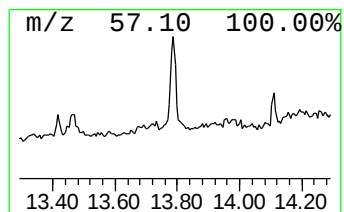
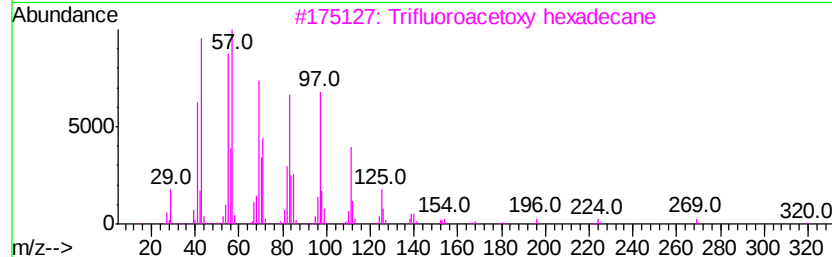
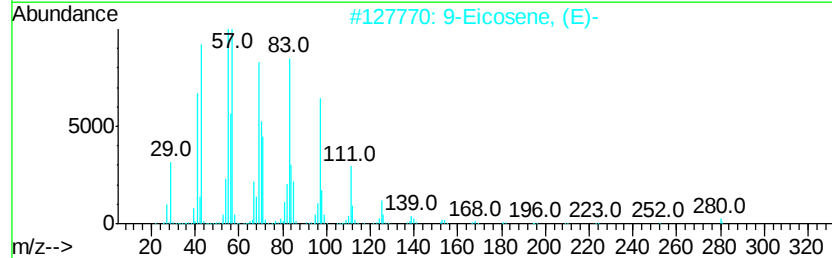
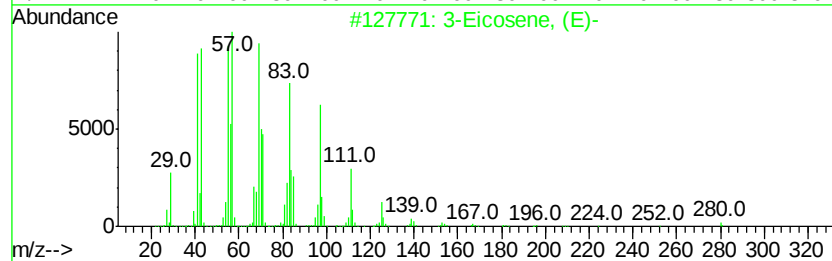
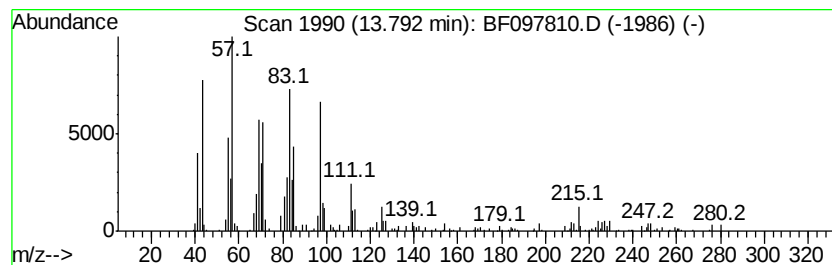
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 3-Eicosene, (E)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.79	4.95 ng	234205	Chrysene-d12	13.93	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-	280	C20H40	074685-33-9	97
2	9-Eicosene, (E)-	280	C20H40	074685-29-3	91
3	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	91
4	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	91
5	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081717\
Data File : BF097810.D
Acq On : 18 Aug 2017 2:34
Operator : SJ/JU
Sample : I4774-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

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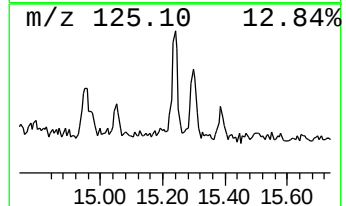
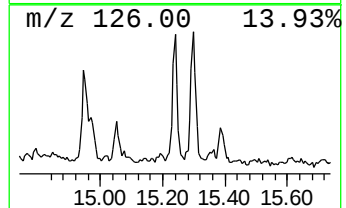
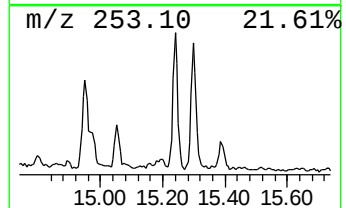
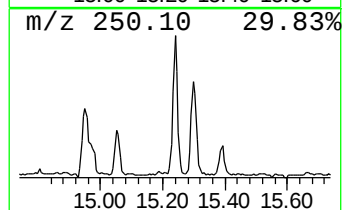
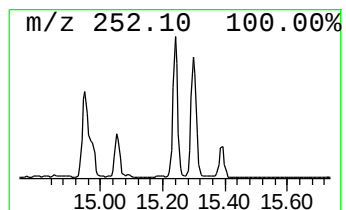
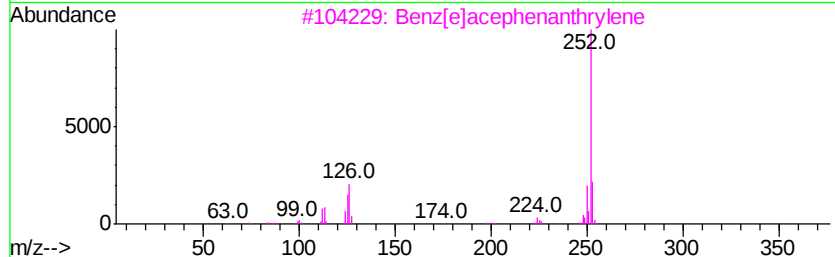
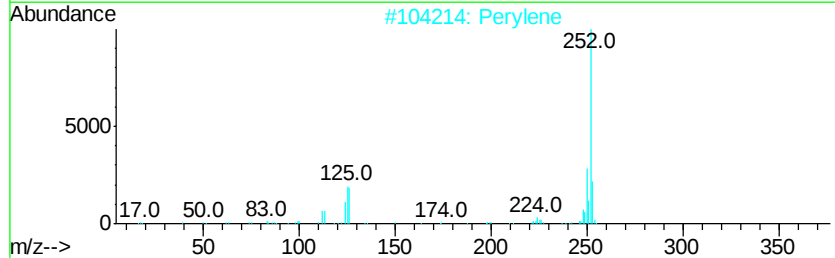
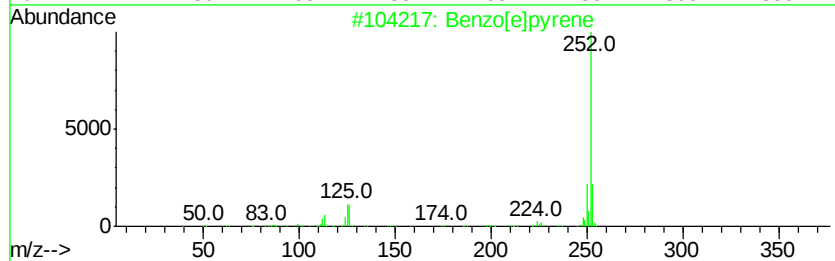
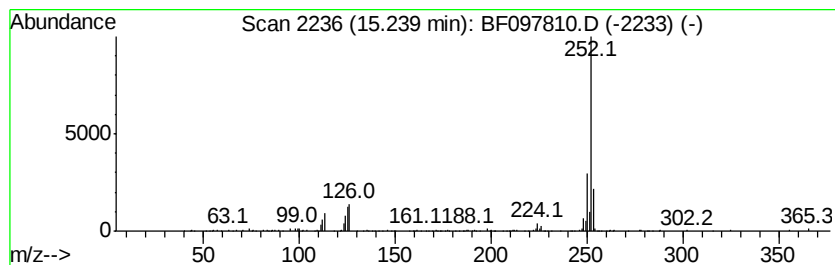
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.24	9.58 ng	215932	Perylene-d12	15.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	97
2			Perylene	252	C20H12	000198-55-0	95
3			Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	91
4			Benzo[a]pyrene	252	C20H12	000050-32-8	91
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	87



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Acq On : 18 Aug 2017 2:34
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Misc :
ALS Vial : 26 Sample Multiplier: 1

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown6.55	6.55	81.8	ng	3356810	1	6.77	820659	20.0
Benzene, 1-etheny...	6.61	2.9	ng	118400	1	6.77	820659	20.0
Naphthalene, 1,7-...	9.40	3.1	ng	131529	3	9.81	833873	20.0
Naphthalene, 2,7-...	9.47	2.3	ng	97235	3	9.81	833873	20.0
Tridecane, 6-propyl-	10.25	2.2	ng	92503	3	9.81	833873	20.0
Tridecane	10.72	6.5	ng	222643	4	11.30	686647	20.0
Phenanthrene, 2-m...	11.83	3.7	ng	126215	4	11.30	686647	20.0
3-Eicosene, (E)-	13.79	5.0	ng	234205	5	13.93	945644	20.0
Benzo[e]pyrene	15.24	9.6	ng	215932	6	15.36	450570	20.0