

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070216\
 Data File : BF088620.D
 Acq On : 2 Jul 2016 16:40
 Operator : UM/SJ
 Sample : H3871-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 L8-B1(0-12)C

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-BF063016.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	1.724	10	12	20	rVB	24118	54534	1.96%	0.193%
2	1.861	22	24	33	rVV	241063	420319	15.09%	1.487%
3	1.998	35	36	46	rVB2	13528	28989	1.04%	0.103%
4	2.993	120	123	132	rBV	26271	58154	2.09%	0.206%
5	4.273	233	235	239	rBV	110233	154855	5.56%	0.548%
6	4.959	292	295	299	rBV	229710	308133	11.06%	1.090%
7	5.382	329	332	335	rBV	1875679	2500786	89.77%	8.845%
8	6.433	420	424	426	rBV	1978535	2785751	100.00%	9.853%
9	6.559	432	435	437	rBV	2606125	2428910	87.19%	8.591%
10	6.719	447	449	451	rBB	32862	29720	1.07%	0.105%
11	6.787	453	455	457	rBV	572320	649793	23.33%	2.298%
12	6.947	467	469	471	rVB	2111929	1915621	68.76%	6.776%
13	7.359	502	505	507	rBV	1818329	1835944	65.90%	6.494%
14	8.079	565	568	570	rBV	1099847	932648	33.48%	3.299%
15	9.153	660	662	665	rVB	2460169	2424653	87.04%	8.576%
16	9.553	695	697	699	rBV	279136	323487	11.61%	1.144%
17	9.771	715	716	719	rBV	45182	40367	1.45%	0.143%
18	9.828	719	721	723	rVV	1101420	920861	33.06%	3.257%
19	9.862	723	724	726	rVB	70647	51950	1.86%	0.184%
20	10.033	737	739	740	rVB	41275	33520	1.20%	0.119%
21	10.273	758	760	761	rVB	61416	48100	1.73%	0.170%
22	10.376	767	769	772	rVB	33434	44388	1.59%	0.157%
23	10.456	775	776	778	rBV2	29083	42880	1.54%	0.152%
24	10.616	787	790	792	rBV	1207834	1387875	49.82%	4.909%
25	10.662	792	794	796	rVB	92159	127059	4.56%	0.449%
26	10.742	800	801	805	rVB	72756	78511	2.82%	0.278%
27	10.891	811	814	816	rBV2	17642	29489	1.06%	0.104%
28	11.188	839	840	846	rVB	102598	145286	5.22%	0.514%
29	11.302	848	850	854	rVV2	734938	1126772	40.45%	3.985%
30	11.382	854	857	858	rVB	122903	140394	5.04%	0.497%
31	11.542	869	871	875	rVB	70903	108227	3.89%	0.383%
32	11.611	875	877	881	rVB2	42270	64492	2.32%	0.228%
33	11.805	892	894	895	rBV	56285	45904	1.65%	0.162%
34	11.839	895	897	899	rVV2	90810	143645	5.16%	0.508%

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35	11.874	899	900	902	rVV	27494	33267	1.19%	0.118%
36	11.919	902	904	908	rVV2	77942	125260	4.50%	0.443%
37	12.022	908	913	915	rVV2	19698	38927	1.40%	0.138%
38	12.102	918	920	925	rVB3	50531	79460	2.85%	0.281%
39	12.354	940	942	944	rBV	32500	37361	1.34%	0.132%
40	12.434	948	949	952	rVB	25690	28131	1.01%	0.099%
41	12.514	954	956	958	rVB	559215	503260	18.07%	1.780%
42	12.571	958	961	964	rBV4	12414	29910	1.07%	0.106%
43	12.628	964	966	968	rBV2	69808	79439	2.85%	0.281%
44	12.742	973	976	978	rBV	325755	517246	18.57%	1.829%
45	12.788	978	980	983	rVB3	24479	45254	1.62%	0.160%
46	12.891	987	989	991	rBV	1863482	1647127	59.13%	5.826%
47	12.959	991	995	999	rVB3	23761	36175	1.30%	0.128%
48	13.062	1000	1004	1006	rVB2	52900	94038	3.38%	0.333%
49	13.131	1009	1010	1012	rVB	63788	54462	1.96%	0.193%
50	13.165	1012	1013	1015	rBV	50333	61548	2.21%	0.218%
51	13.257	1019	1021	1023	rBV	25385	30338	1.09%	0.107%
52	13.291	1023	1024	1026	rBV	33271	30086	1.08%	0.106%
53	13.474	1038	1040	1042	rBV2	27566	31664	1.14%	0.112%
54	13.565	1046	1048	1051	rVB	33251	46496	1.67%	0.164%
55	13.679	1055	1058	1060	rBV2	43246	74263	2.67%	0.263%
56	13.794	1067	1068	1072	rVB2	95016	136626	4.90%	0.483%
57	13.931	1077	1080	1086	rVB2	812653	1100817	39.52%	3.894%
58	14.034	1087	1089	1091	rBV2	38109	37416	1.34%	0.132%
59	14.114	1095	1096	1098	rBV	35979	51340	1.84%	0.182%
60	14.320	1112	1114	1115	rVB	32309	35304	1.27%	0.125%
61	14.788	1153	1155	1157	rVB	71261	72224	2.59%	0.255%
62	14.937	1166	1168	1173	rVB	169739	345263	12.39%	1.221%
63	15.211	1190	1192	1195	rVB	114650	150868	5.42%	0.534%
64	15.268	1195	1197	1199	rBV	145257	178052	6.39%	0.630%
65	15.325	1200	1202	1204	rBV	363914	516146	18.53%	1.826%
66	16.000	1258	1261	1265	rVB	76831	169771	6.09%	0.600%
67	16.400	1293	1296	1300	rVB	75838	157814	5.67%	0.558%
68	16.663	1317	1319	1323	rVB2	55042	102820	3.69%	0.364%
69	16.926	1340	1342	1345	rVB	45619	78625	2.82%	0.278%
70	17.063	1352	1354	1359	rVB	39721	83862	3.01%	0.297%

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Peak separation: 5

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

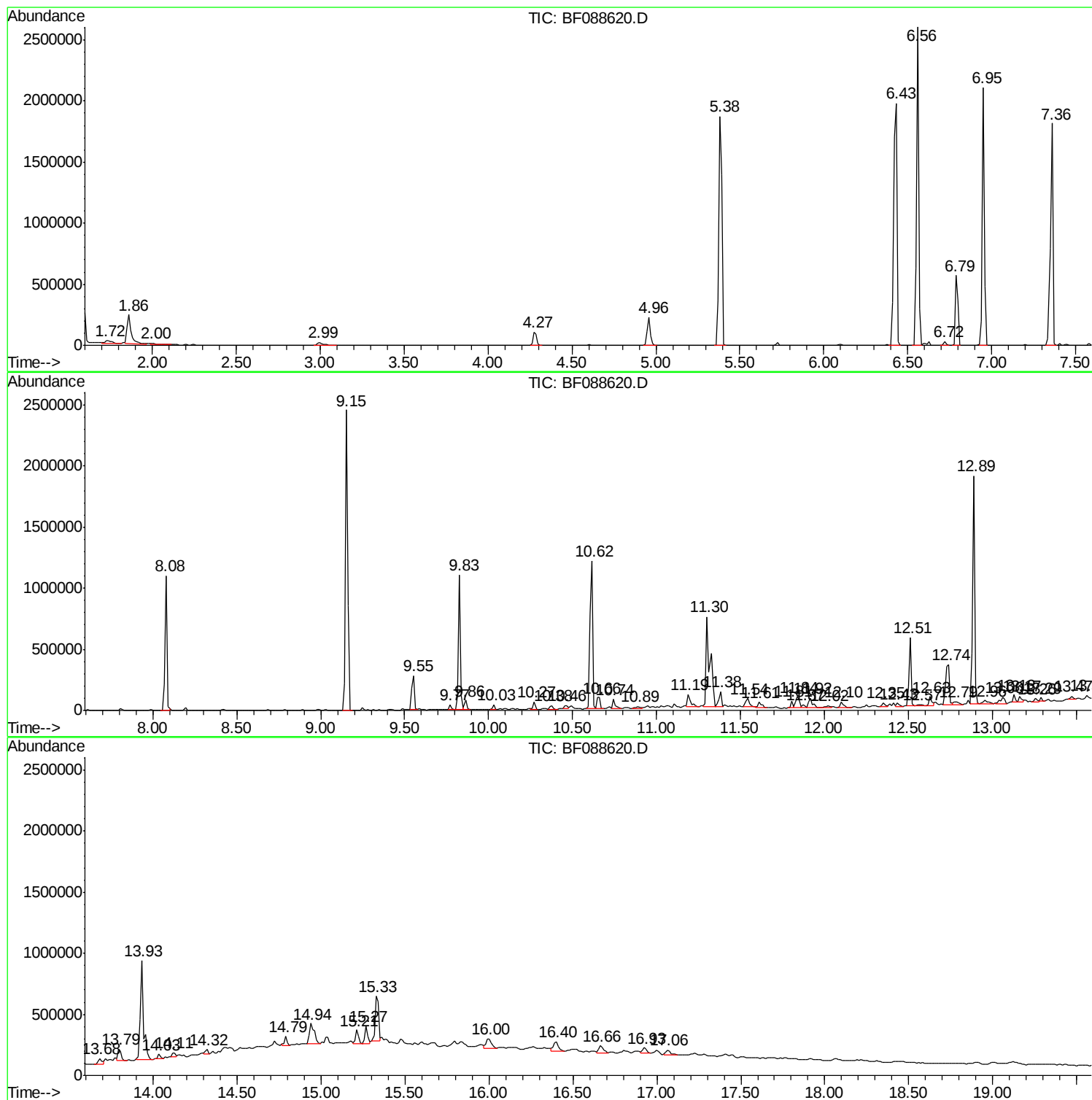
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Sample : H3871-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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L8-B1(0-12)C

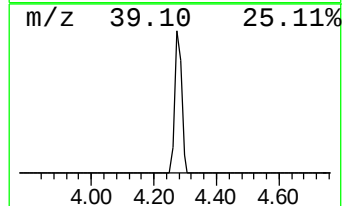
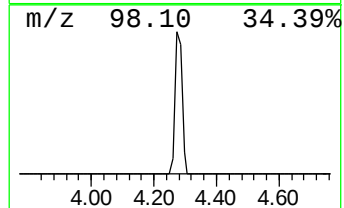
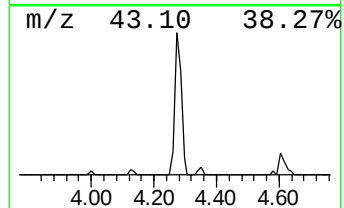
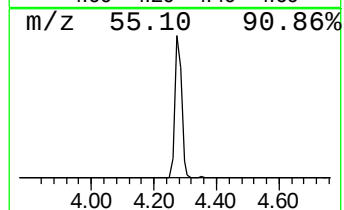
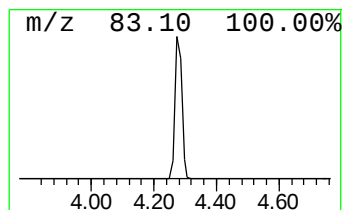
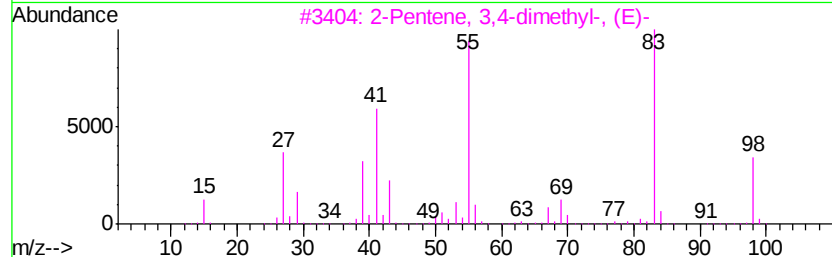
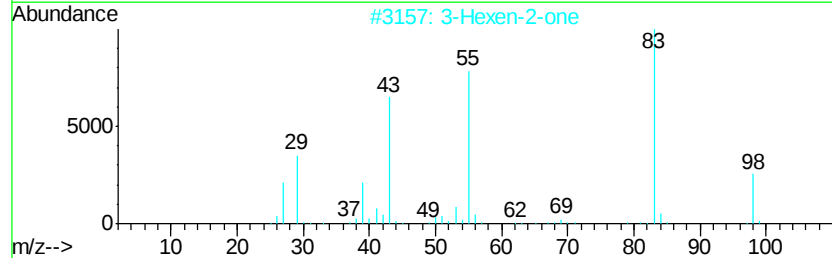
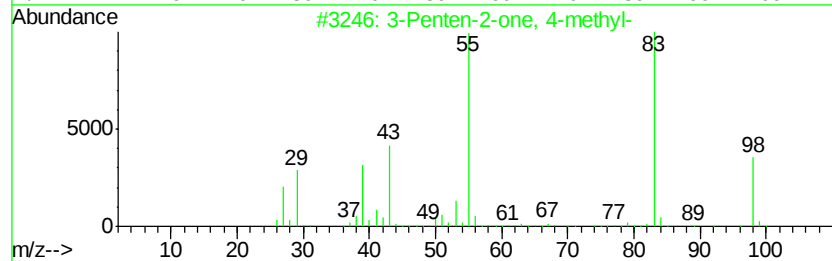
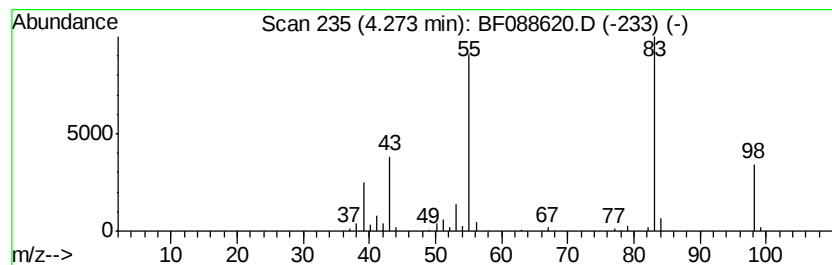
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.27	4.77 ng	154855	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	95
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, 2,4-dimethyl-	98	C7H14	000625-65-0	87
5			2-Pentene, 4,4-dimethyl-, (Z)-	98	C7H14	000762-63-0	86



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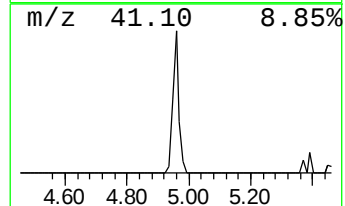
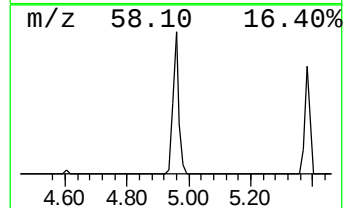
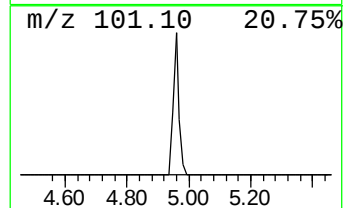
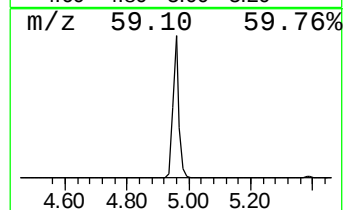
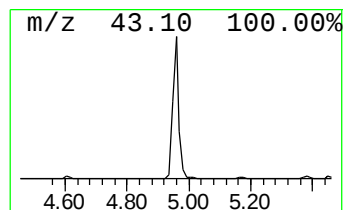
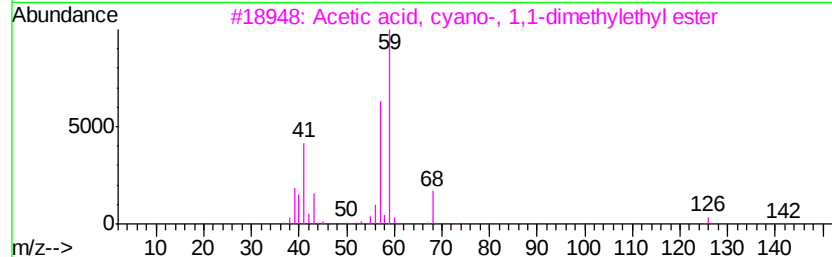
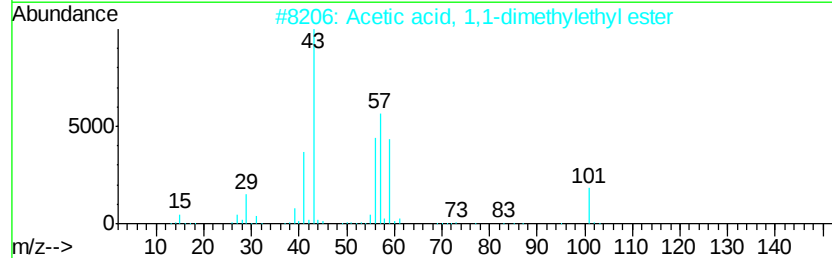
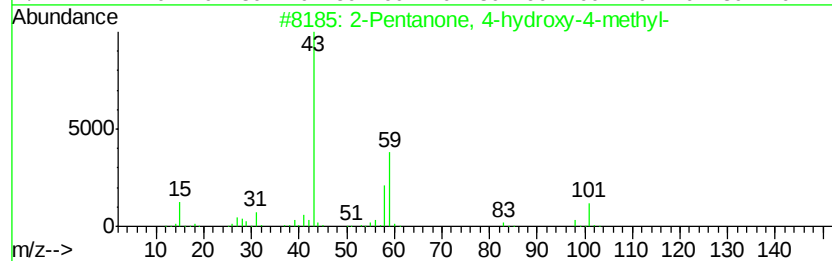
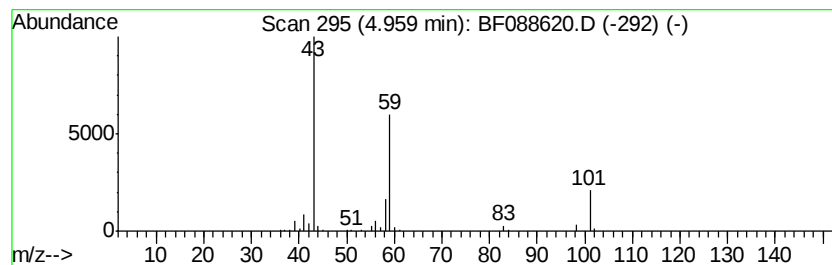
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	9.48 ng	308133	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
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	32
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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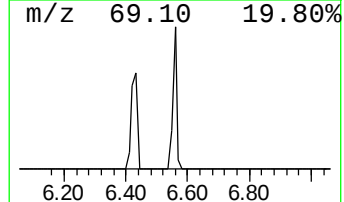
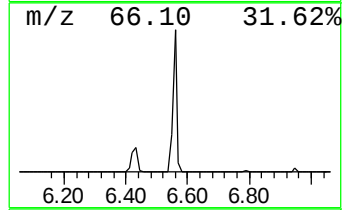
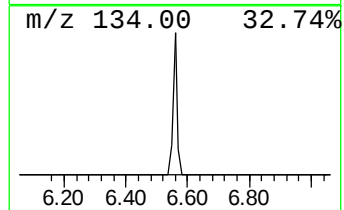
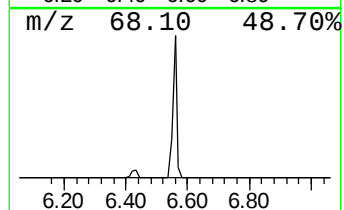
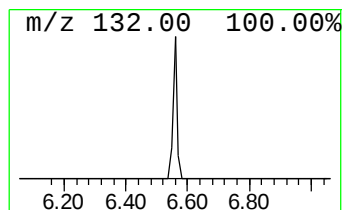
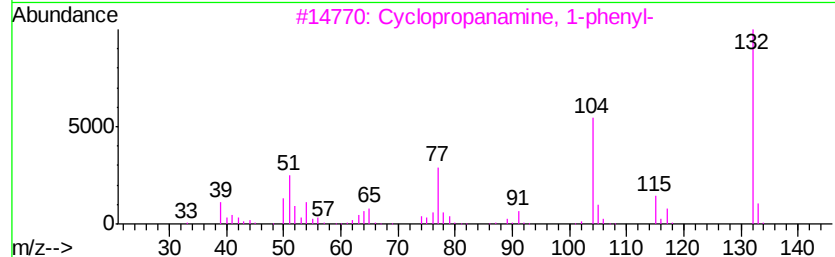
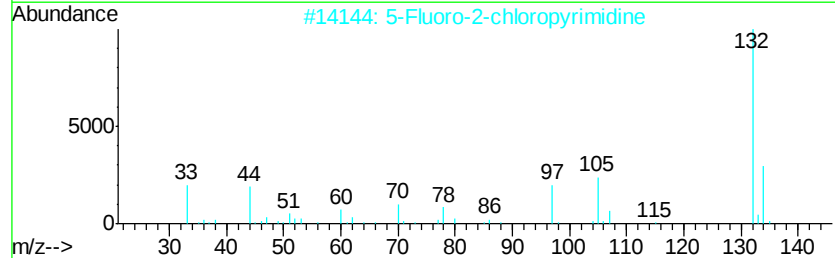
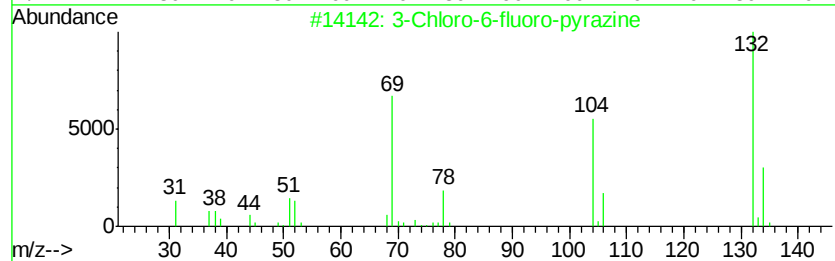
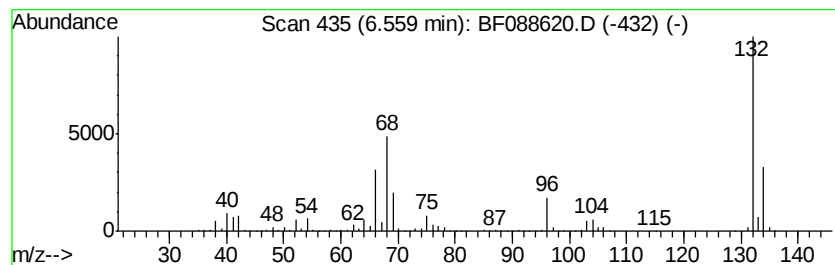
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	74.76 ng	2428910	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	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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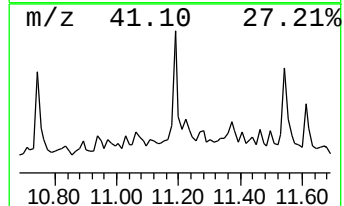
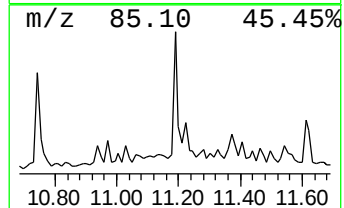
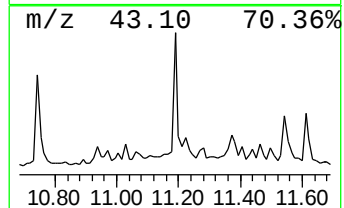
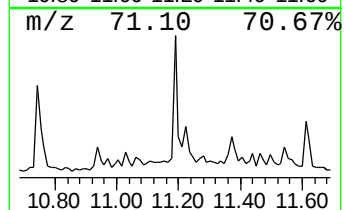
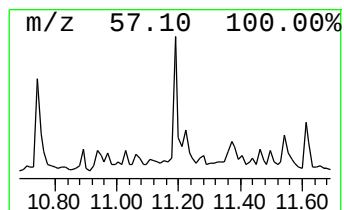
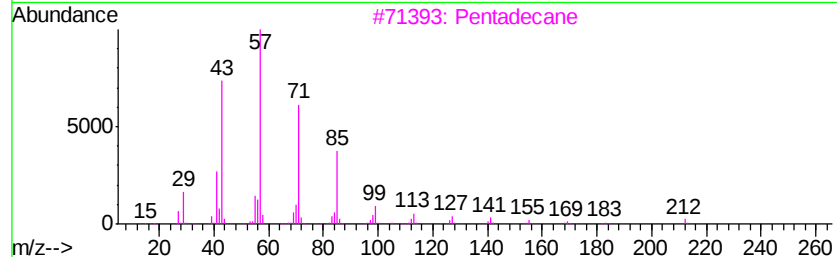
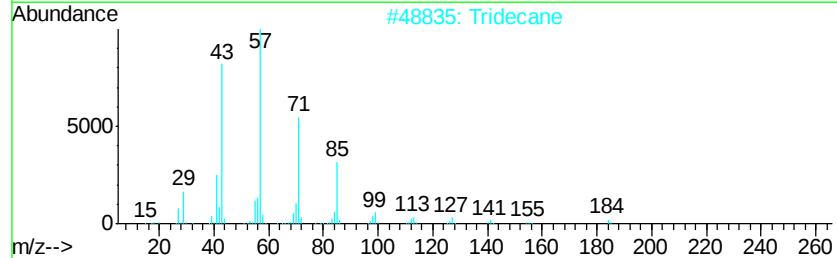
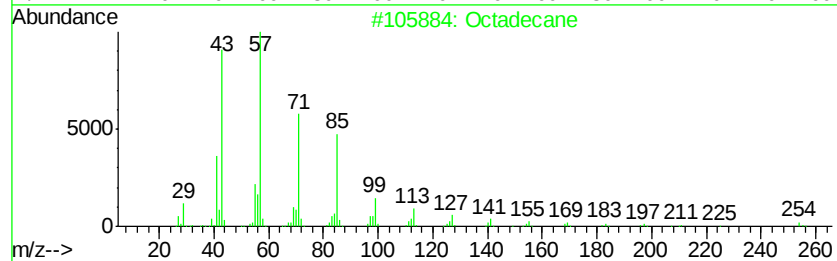
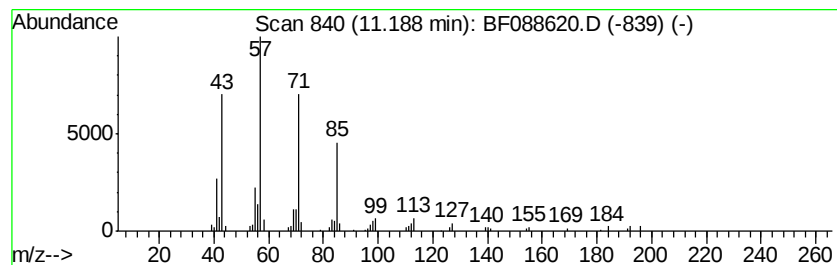
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L8-B1(0-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Octadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.		
11.19	2.58 ng	145286	Phenanthrene-d10		11.30		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	86
2			Tridecane	184	C13H28	000629-50-5	81
3			Pentadecane	212	C15H32	000629-62-9	80
4			Heneicosane	296	C21H44	000629-94-7	78
5			Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070216\
Data File : BF088620.D
Acq On : 2 Jul 2016 16:40
Operator : UM/SJ
Sample : H3871-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
L8-B1(0-12)C

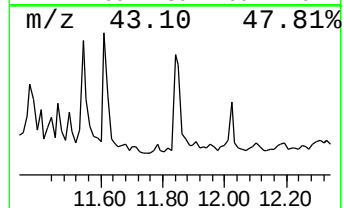
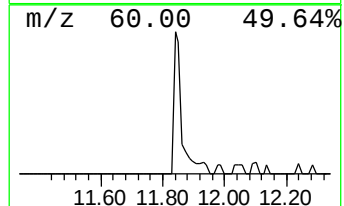
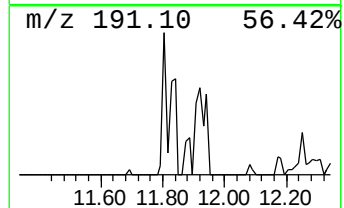
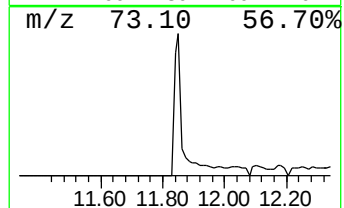
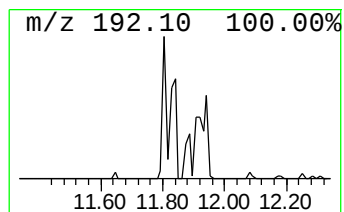
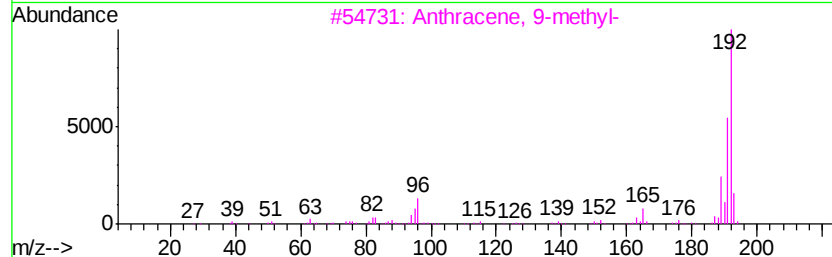
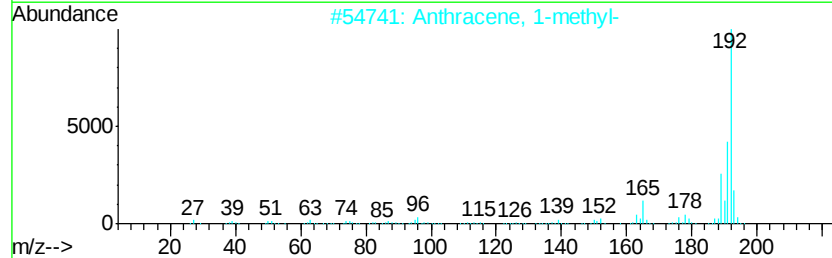
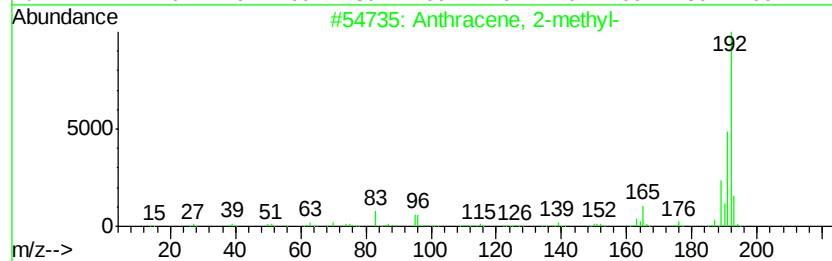
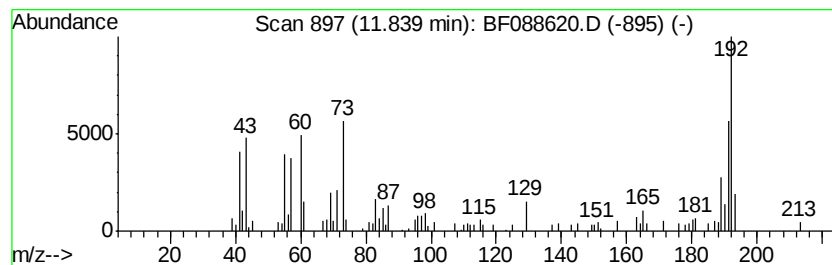
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Anthracene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.84	2.55 ng	143645	Phenanthrene-d10	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
3		Anthracene, 9-methyl-	192	C15H12	000779-02-2	83
4		1H-Cyclopropa[1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	78
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070216\
Data File : BF088620.D
Acq On : 2 Jul 2016 16:40
Operator : UM/SJ
Sample : H3871-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

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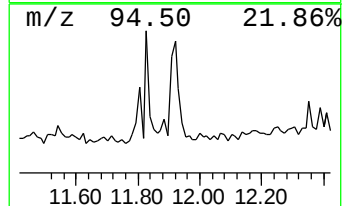
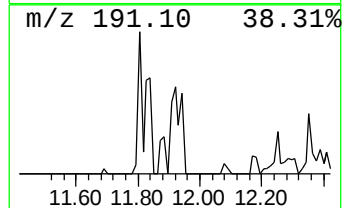
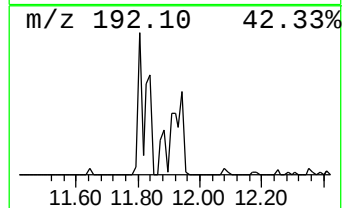
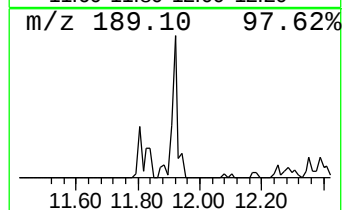
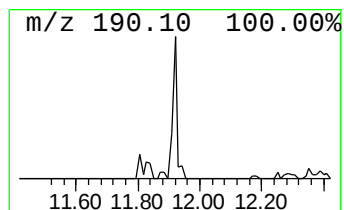
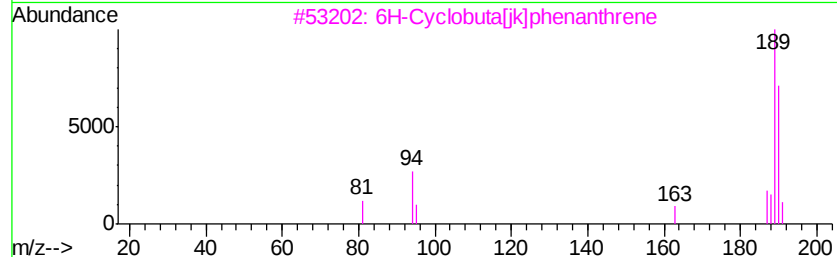
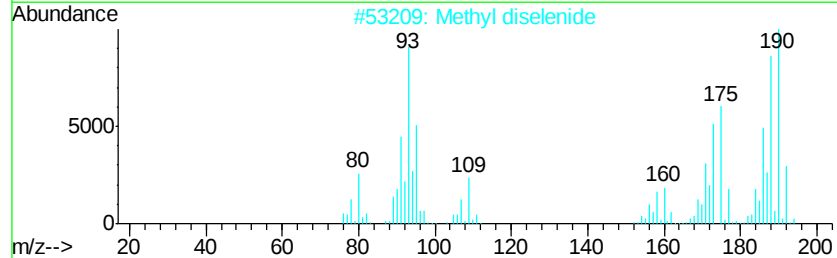
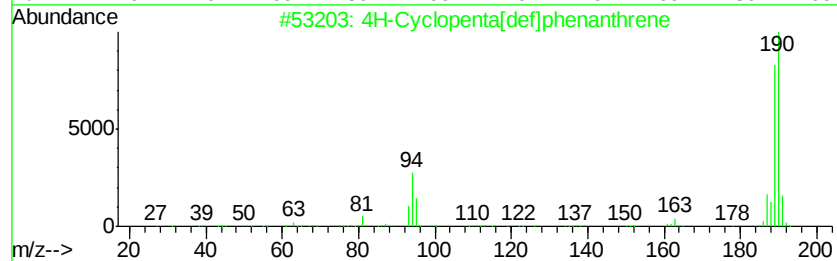
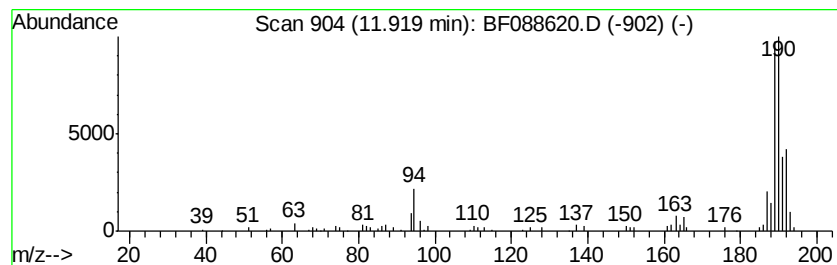
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	2.22 ng	125260	Phenanthrene-d10	11.30

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	62
2		Methyl diselenide	190	C2H6Se2	007101-31-7	45
3		6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	42
4		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	37
5		2,6-Dichlorobenzaldoxime	189	C7H5Cl2NO	025185-95-9	30



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Operator : UM/SJ
Sample : H3871-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
L8-B1(0-12)C

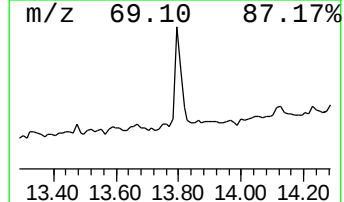
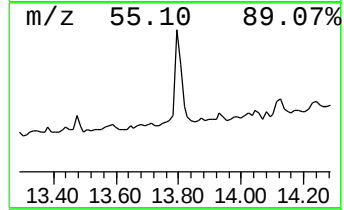
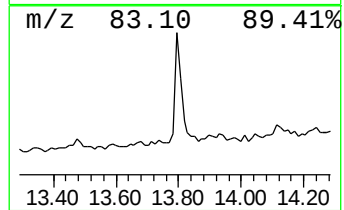
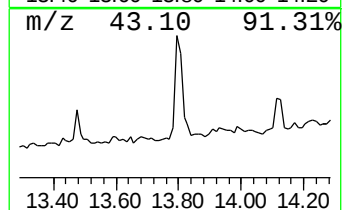
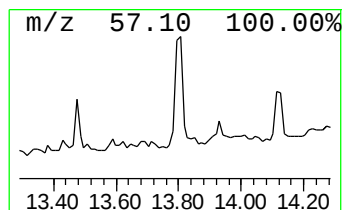
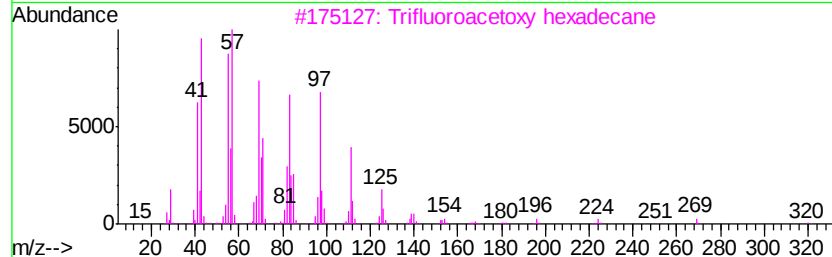
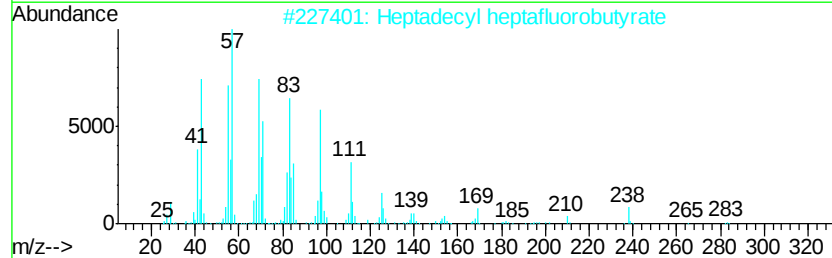
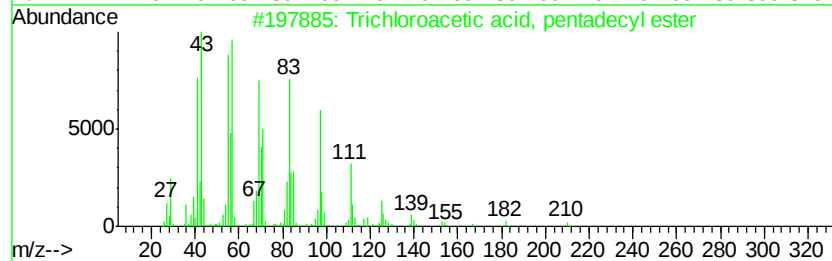
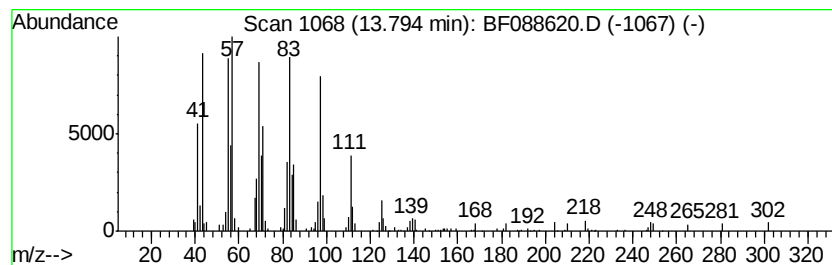
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Trichloroacetic acid, penta... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	2.48 ng	136626	Chrysene-d12	13.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	94
2		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
4		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	94
5		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94



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Sample : H3871-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
L8-B1(0-12)C

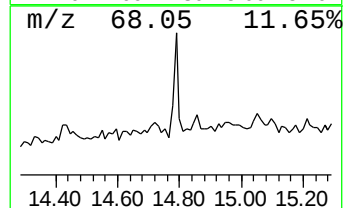
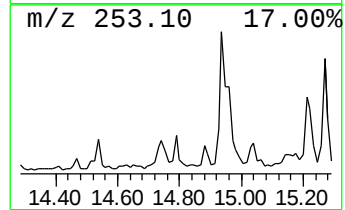
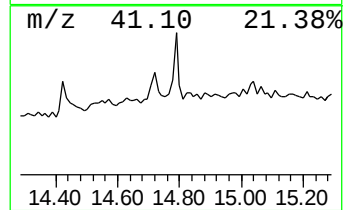
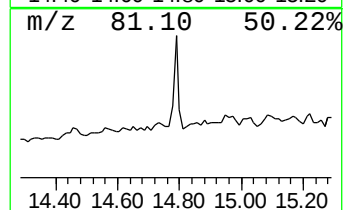
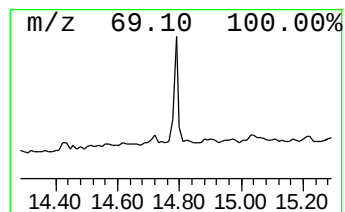
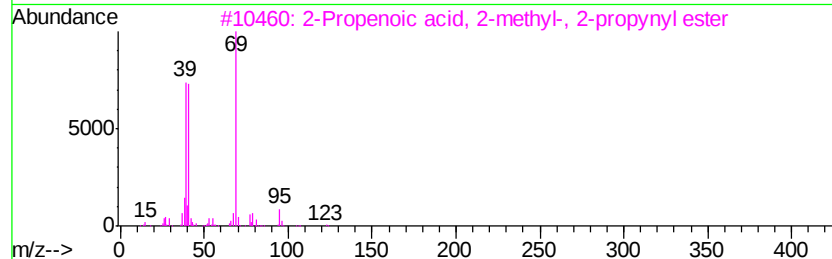
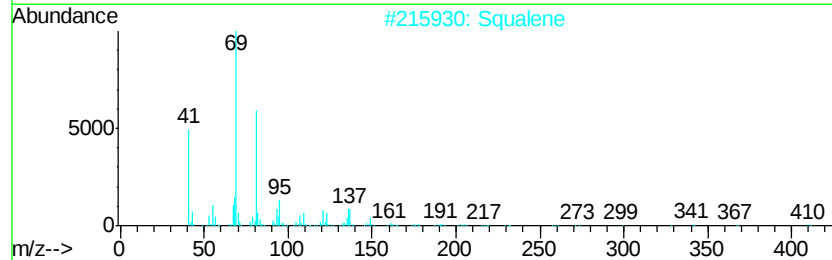
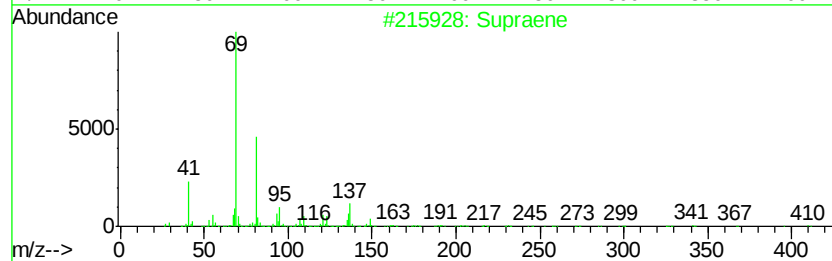
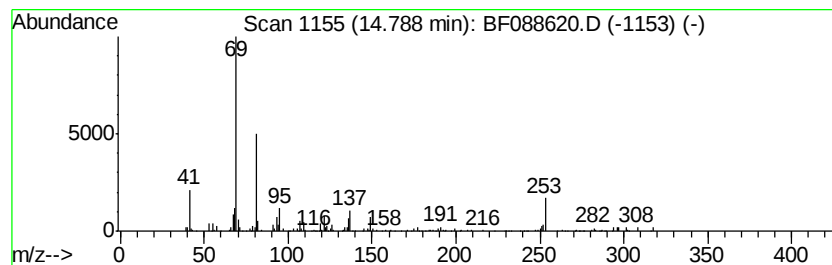
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Supraene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.79	2.80 ng	72224	Perylene-d12	15.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	74
2			Squalene	410	C30H50	000111-02-4	72
3			2-Propenoic acid, 2-methyl-, 2-p...	124	C7H8O2	013861-22-8	53
4			4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	50
5			3,7,11-Tridecatrienitrile, 4,8...	231	C16H25N	006006-01-5	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070216\
Data File : BF088620.D
Acq On : 2 Jul 2016 16:40
Operator : UM/SJ
Sample : H3871-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
L8-B1(0-12)C

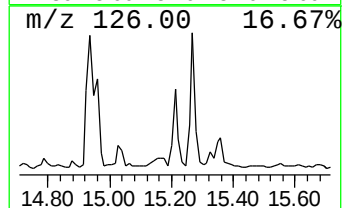
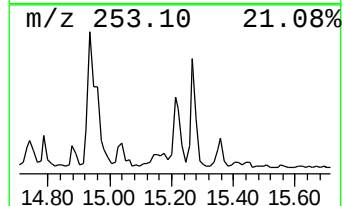
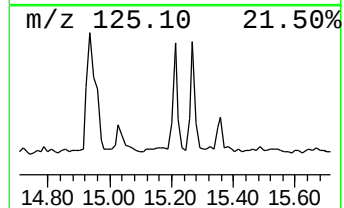
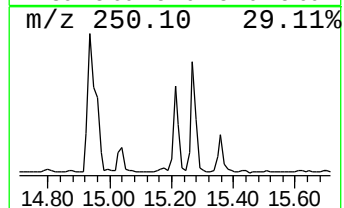
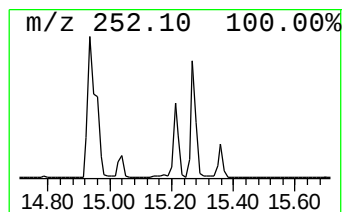
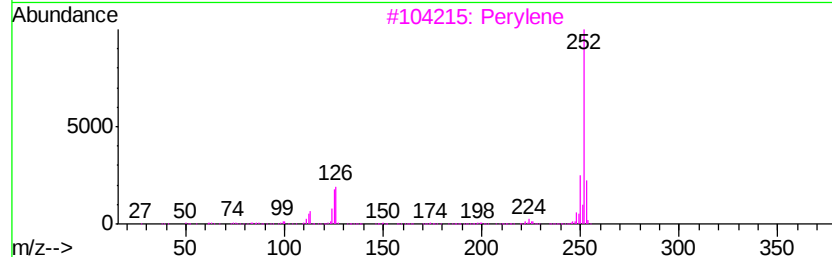
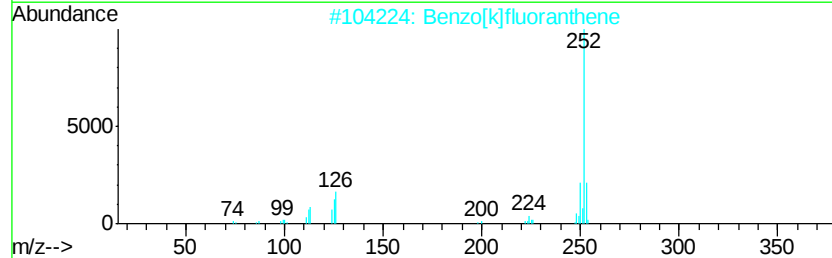
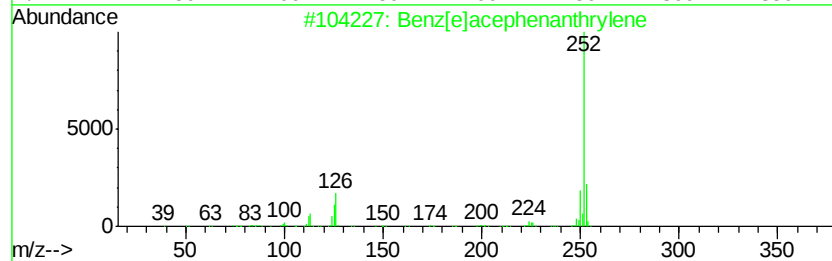
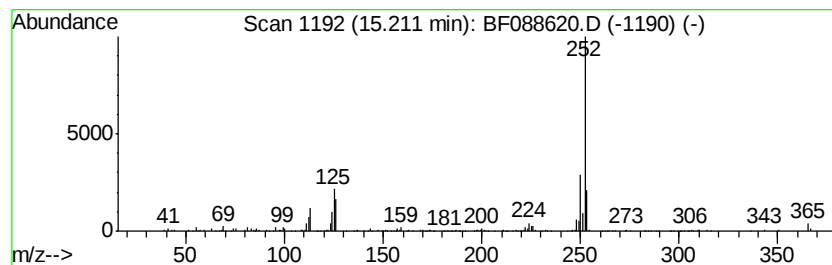
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benz[e]acephenanthrylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.21	5.85 ng	150868	Perylene-d12	15.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	96
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
3			Perylene	252	C20H12	000198-55-0	93
4			Benzo[e]pyrene	252	C20H12	000192-97-2	93
5			Benzo[j]fluoranthene	252	C20H12	000205-82-3	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070216\
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Misc :
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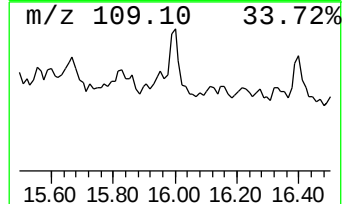
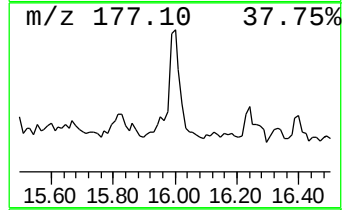
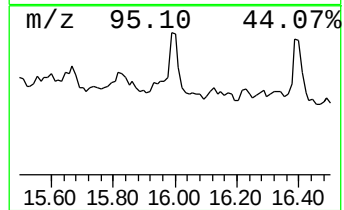
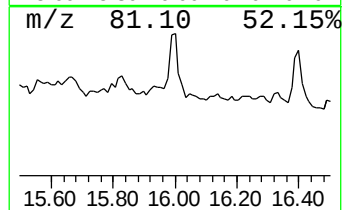
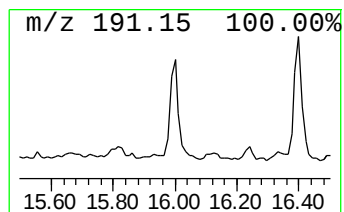
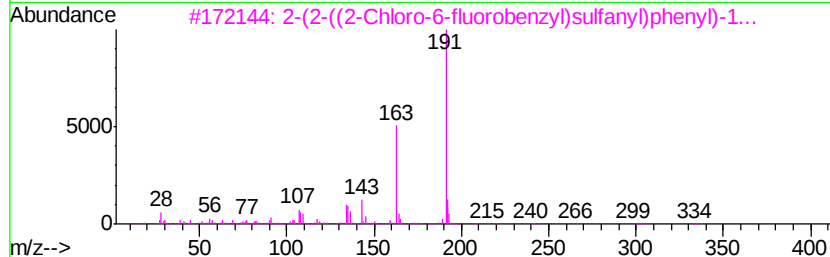
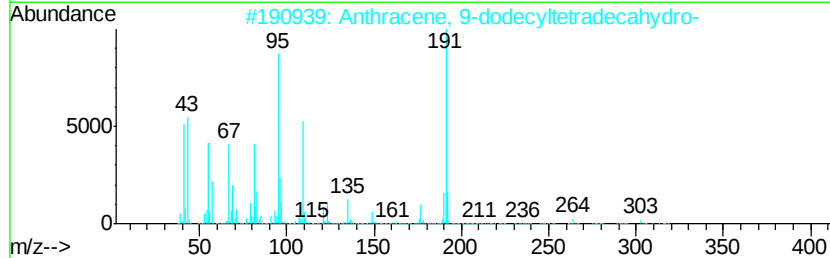
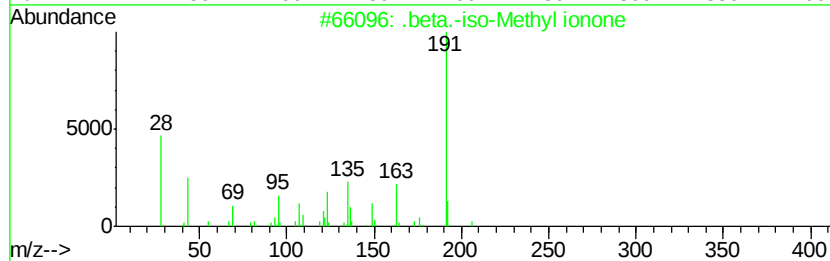
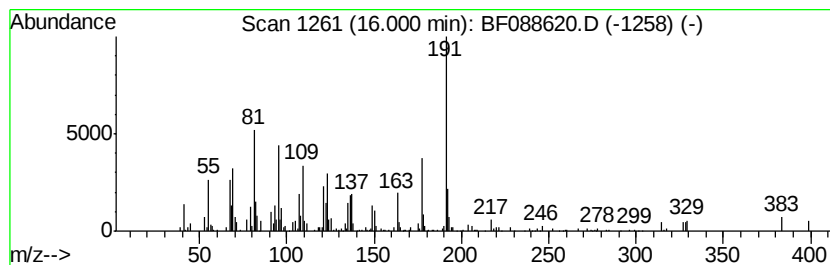
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 .beta.-iso-Methyl ionone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.00	6.58 ng	169771	Perylene-d12	15.34	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	64
2	Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	38
3	2-(2-((2-Chloro-6-fluorobenzyl)s...	334	C17H16ClFN2S	1000298-18-2	38
4	1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	37
5	Pentanoic acid, 4-methyl-2-(2,5-...	278	C13H18N4O3	1000294-63-5	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070216\
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ALS Vial : 11 Sample Multiplier: 1

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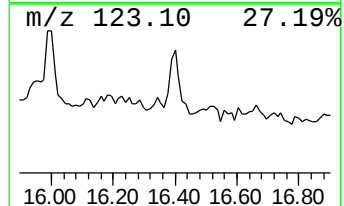
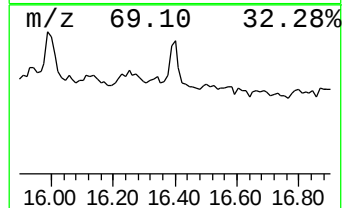
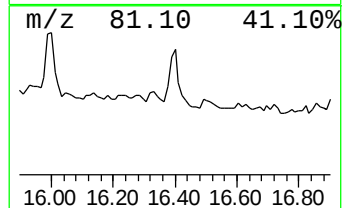
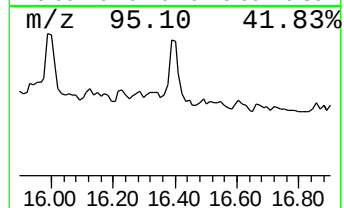
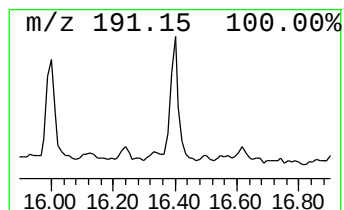
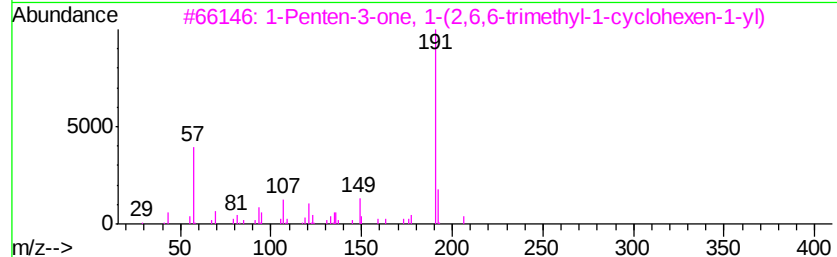
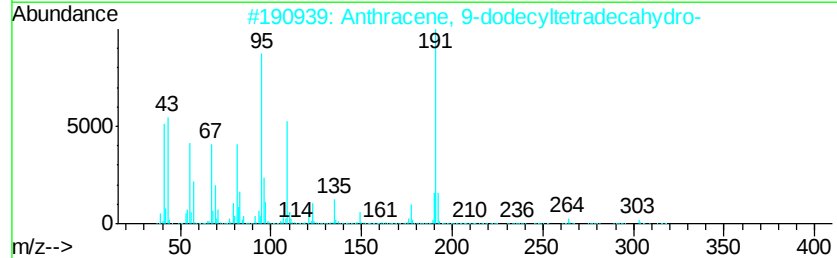
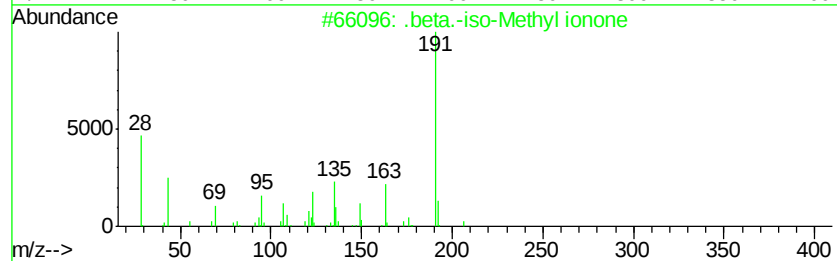
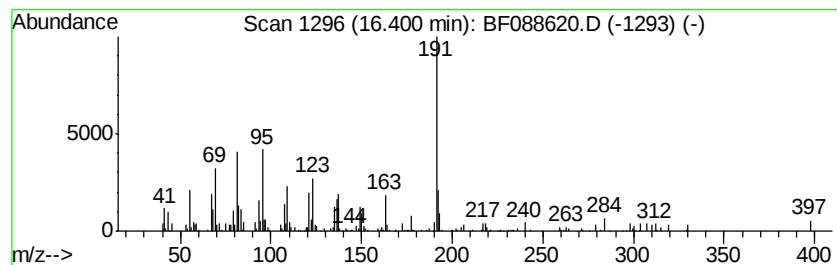
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown16.40 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.40	6.12 ng	157814	Perylene-d12	15.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	43
2		Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	42
3		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	38
4		9H-Fluorene-9-carbonitrile	191	C14H9N	001529-40-4	35
5		Phenol, 2,4-bis(1,1-dimethylethyl)-	206	C14H22O	000096-76-4	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070216\
Data File : BF088620.D
Acq On : 2 Jul 2016 16:40
Operator : UM/SJ
Sample : H3871-09
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
L8-B1(0-12)C

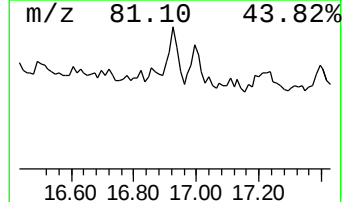
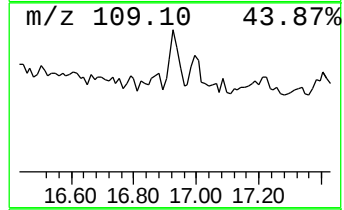
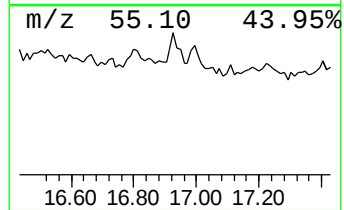
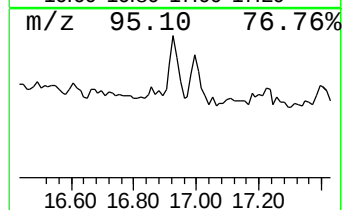
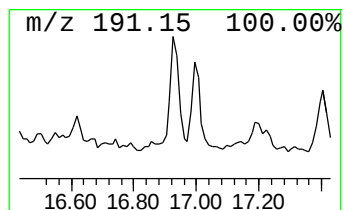
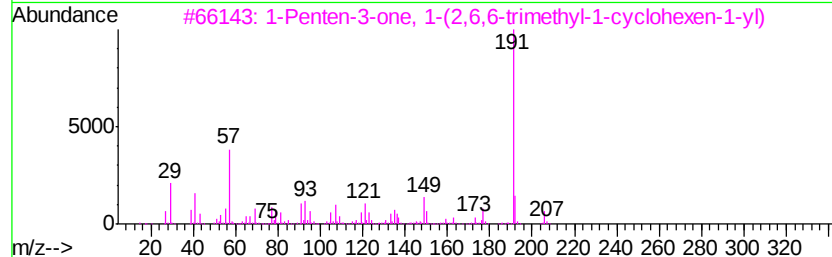
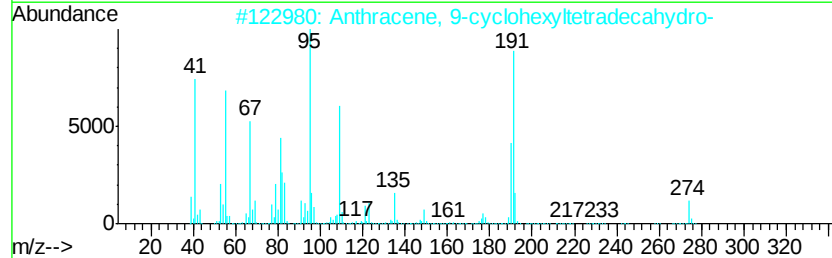
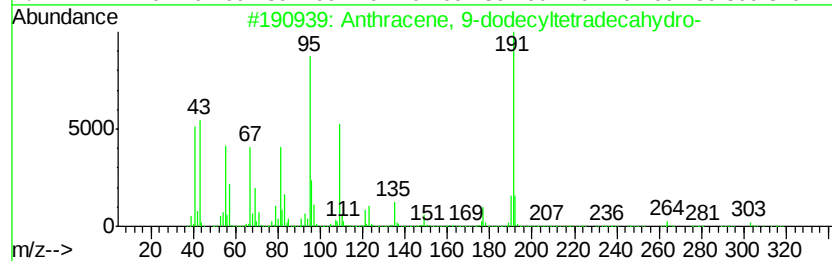
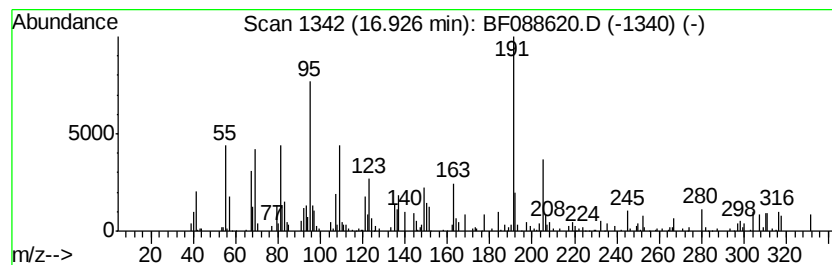
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Anthracene, 9-dodecyltetrad... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.93	3.05 ng	78625	Perylene-d12	15.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	64
2			Anthracene, 9-cyclohexyltetradec...	274	C20H34	055255-70-4	53
3			1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	30
4			4H-Benzo[def]carbazole	191	C14H9N	000203-65-6	27
5			Iridomyrmecin	168	C10H16O2	000485-43-8	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070216\
 Data File : BF088620.D
 Acq On : 2 Jul 2016 16:40
 Operator : UM/SJ
 Sample : H3871-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 L8-B1(0-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.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...	4.27	4.8 ng		154855	1	6.79	649793	20.0
2-Pentanone, 4-hy...	4.96	9.5 ng		308133	1	6.79	649793	20.0
unknown6.56	6.56	74.8 ng		2428910	1	6.79	649793	20.0
Octadecane	11.19	2.6 ng		145286	4	11.30	1126770	20.0
Anthracene, 2-met...	11.84	2.5 ng		143645	4	11.30	1126770	20.0
4H-Cyclopenta[def...	11.92	2.2 ng		125260	4	11.30	1126770	20.0
Trichloroacetic a...	13.79	2.5 ng		136626	5	13.93	1100820	20.0
Supraene	14.79	2.8 ng		72224	6	15.34	516146	20.0
Benz[e]acephenant...	15.21	5.8 ng		150868	6	15.34	516146	20.0
.beta.-iso-Methyl...	16.00	6.6 ng		169771	6	15.34	516146	20.0
unknown16.40	16.40	6.1 ng		157814	6	15.34	516146	20.0
Anthracene, 9-dod...	16.93	3.0 ng		78625	6	15.34	516146	20.0