

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050516\
 Data File : BF086714.D
 Acq On : 5 May 2016 19:21
 Operator : UM/SJ
 Sample : H2829-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EUT01-TD-B12(0-8.5)-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-BF050416.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	4.887	242	245	252	rBV	317061	489078	9.60%	0.958%
2	5.322	280	283	285	rBV	3925600	5034918	98.83%	9.866%
3	5.527	299	301	309	rVB	84953	115013	2.26%	0.225%
4	5.722	316	318	324	rVB	50669	67835	1.33%	0.133%
5	6.373	372	375	377	rBV	4471383	5055922	99.24%	9.907%
6	6.499	383	386	388	rBV	3820700	5094768	100.00%	9.983%
7	6.556	388	391	394	rVV	95008	153729	3.02%	0.301%
8	6.727	404	406	408	rBV	1111974	1213182	23.81%	2.377%
9	6.876	417	419	422	rBV	2775927	3610026	70.86%	7.074%
10	6.990	427	429	435	rVB2	67265	119633	2.35%	0.234%
11	7.299	452	456	458	rBV	2475365	3349451	65.74%	6.563%
12	7.402	462	465	472	rVB2	96370	176578	3.47%	0.346%
13	7.585	480	481	485	rVB	81517	84515	1.66%	0.166%
14	7.756	494	496	499	rBV	32473	53691	1.05%	0.105%
15	8.007	516	518	521	rBV	1428348	1540004	30.23%	3.018%
16	8.293	541	543	548	rBV	50711	63608	1.25%	0.125%
17	9.093	610	613	615	rBV	4998905	4844663	95.09%	9.493%
18	9.208	622	623	629	rVB	138368	209403	4.11%	0.410%
19	9.482	645	647	650	rBV	484293	498322	9.78%	0.976%
20	9.711	665	667	669	rBV	118964	173317	3.40%	0.340%
21	9.756	669	671	673	rVV	1190740	1650984	32.41%	3.235%
22	9.791	673	674	677	rVB2	89618	100626	1.98%	0.197%
23	10.202	709	710	712	rVB	176979	147926	2.90%	0.290%
24	10.305	718	719	723	rVB	52084	51377	1.01%	0.101%
25	10.419	726	729	732	rBV	66295	109361	2.15%	0.214%
26	10.545	738	740	743	rBV2	2233726	3020151	59.28%	5.918%
27	10.671	750	751	755	rBV	185520	279538	5.49%	0.548%
28	10.865	765	768	772	rBV3	38445	91015	1.79%	0.178%
29	11.128	786	791	796	rVB3	113975	276253	5.42%	0.541%
30	11.242	799	801	802	rBV	1633767	1517895	29.79%	2.974%
31	11.311	805	807	810	rVB2	103232	144877	2.84%	0.284%
32	11.482	819	822	826	rBV3	33308	79149	1.55%	0.155%
33	11.551	826	828	829	rBV2	74282	79325	1.56%	0.155%
34	11.745	843	845	846	rBV	59364	71091	1.40%	0.139%

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EUT01-TD-B12(0-8.5)-C

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35	11.779	846	848	850	rVV3	340057	473789	9.30%	0.928%
36	11.848	853	854	859	rVB	123606	194753	3.82%	0.382%
37	12.042	868	871	876	rVB2	37023	76948	1.51%	0.151%
38	12.294	890	893	894	rBV	44874	58896	1.16%	0.115%
39	12.339	894	897	899	rVB2	28456	58822	1.15%	0.115%
40	12.454	904	907	909	rBV	516579	757572	14.87%	1.484%
41	12.568	915	917	919	rBV2	78762	77369	1.52%	0.152%
42	12.671	923	926	929	rBV	622304	785813	15.42%	1.540%
43	12.728	929	931	935	rVB3	35933	69450	1.36%	0.136%
44	12.831	937	940	942	rBV	3533597	3527567	69.24%	6.912%
45	13.002	952	955	957	rVB2	98446	134057	2.63%	0.263%
46	13.071	959	961	962	rVV	101514	90908	1.78%	0.178%
47	13.105	962	964	968	rVB2	70167	111865	2.20%	0.219%
48	13.414	987	991	996	rBV5	28108	106711	2.09%	0.209%
49	13.619	1007	1009	1010	rBV2	49998	64022	1.26%	0.125%
50	13.734	1017	1019	1022	rBV2	225079	299085	5.87%	0.586%
51	13.871	1028	1031	1037	rBV2	1143098	1677248	32.92%	3.286%
52	13.974	1038	1040	1042	rBV	57392	51474	1.01%	0.101%
53	14.054	1045	1047	1050	rBV4	32970	74029	1.45%	0.145%
54	14.374	1071	1075	1079	rBV5	57222	171516	3.37%	0.336%
55	14.877	1116	1119	1124	rBV	267829	566444	11.12%	1.110%
56	14.968	1124	1127	1132	rVV4	65879	181624	3.56%	0.356%
57	15.140	1140	1142	1145	rBV	133951	190562	3.74%	0.373%
58	15.197	1145	1147	1150	rBV	176456	252842	4.96%	0.495%
59	15.254	1150	1152	1155	rBV	559047	937424	18.40%	1.837%
60	15.688	1188	1190	1192	rBV	59794	97481	1.91%	0.191%
61	16.568	1264	1267	1272	rVB2	95172	197194	3.87%	0.386%
62	16.957	1298	1301	1308	rBV	83822	182195	3.58%	0.357%

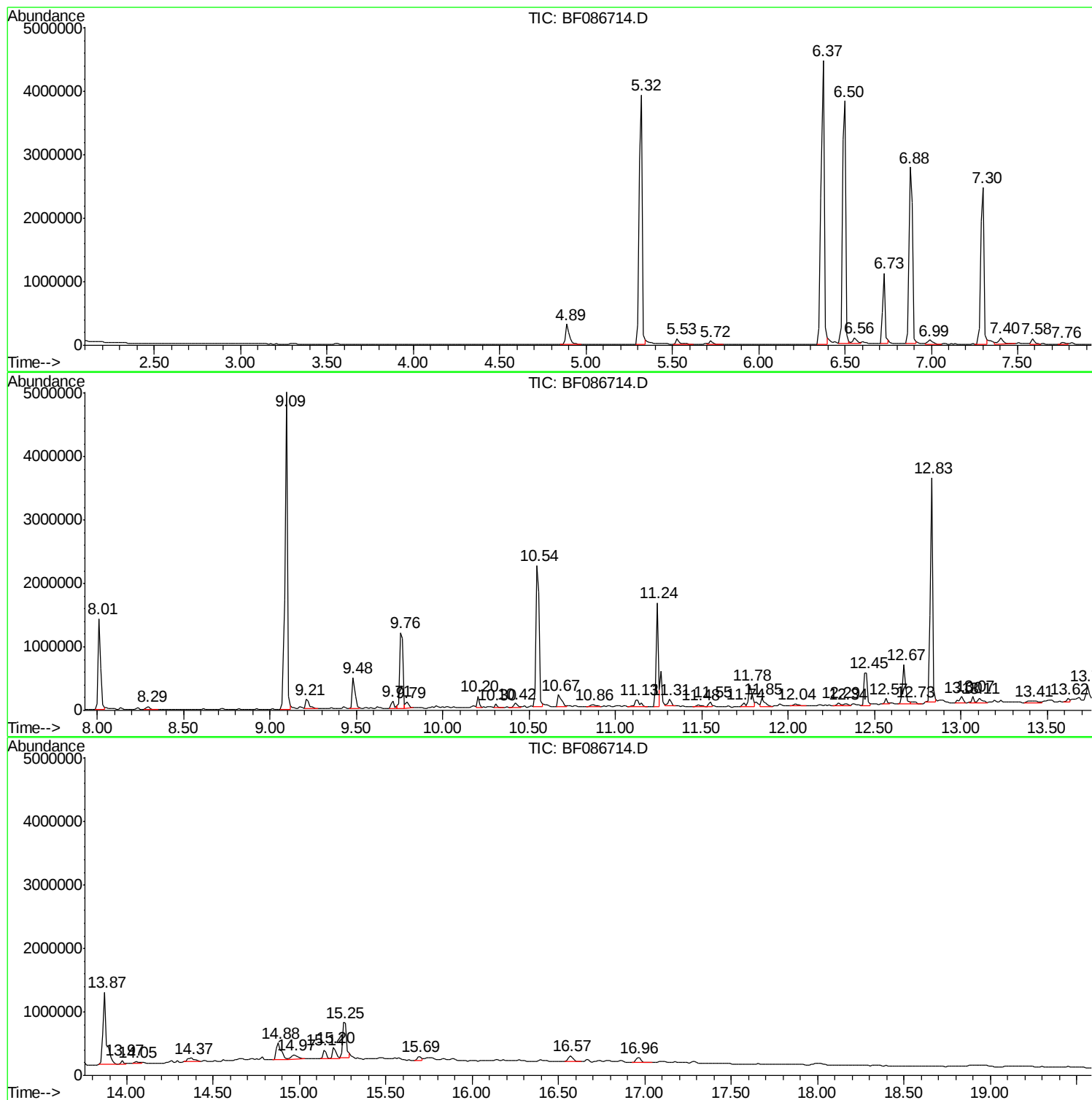
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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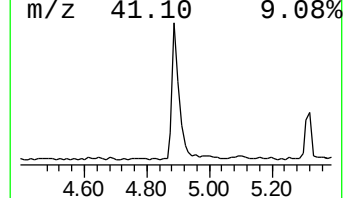
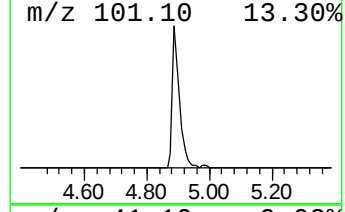
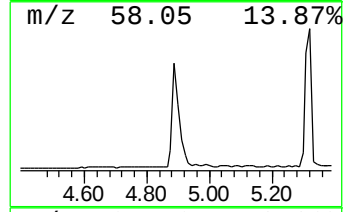
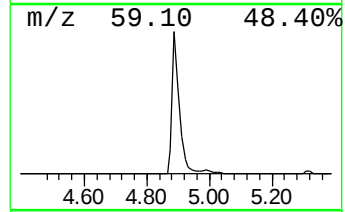
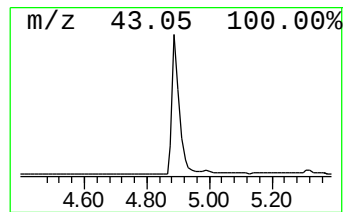
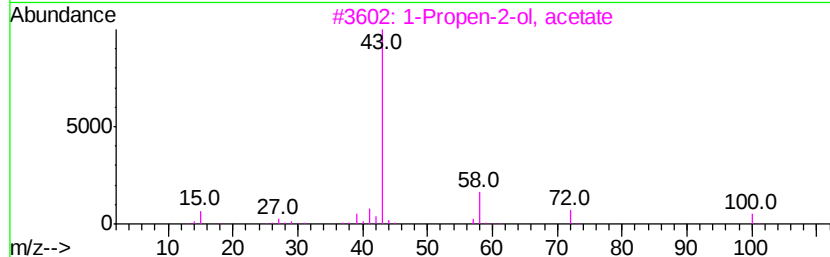
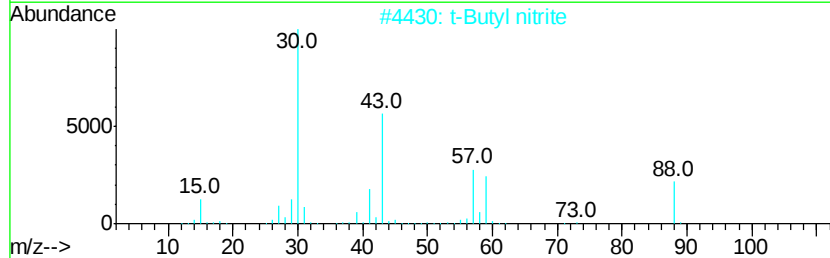
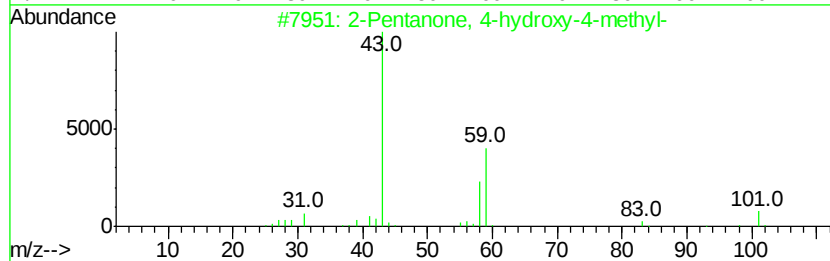
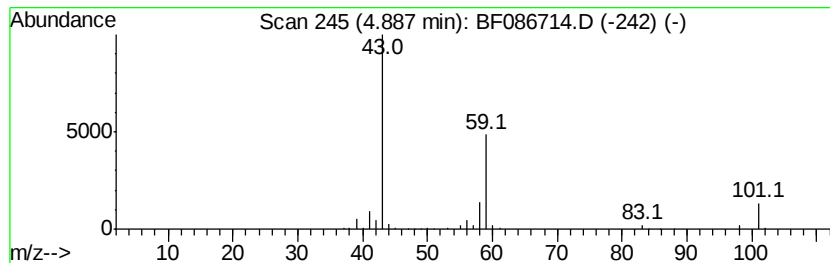
Instrument :
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ClientSampleId :
EUT01-TD-B12(0-8.5)-C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
4.89	16.13 ng	489078	1,4-Dichlorobenzene-d4	6.73		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50	
2	t-Butyl nitrite	103	C4H9NO2	000540-80-7	39	
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10	
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9	
5	2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9	



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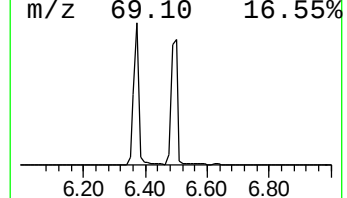
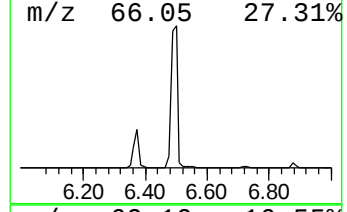
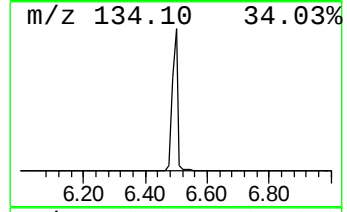
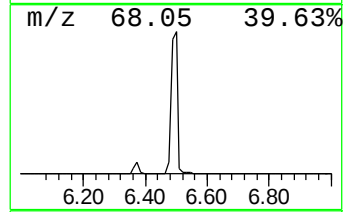
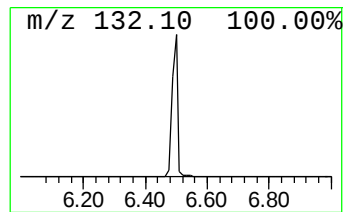
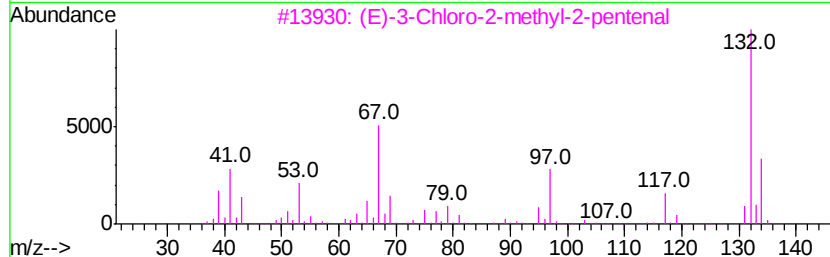
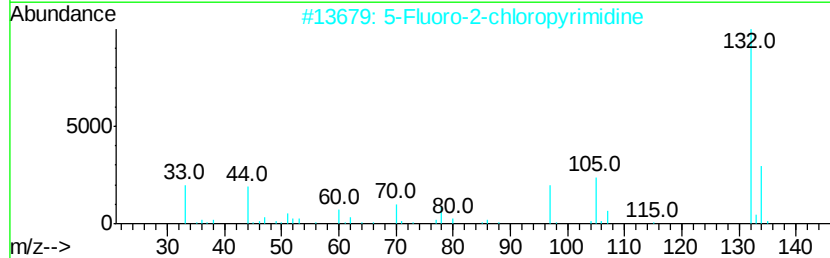
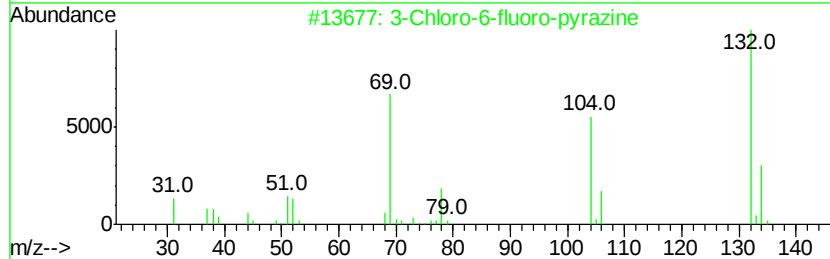
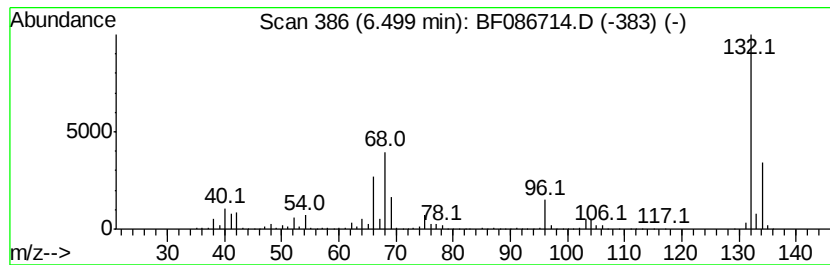
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.50 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	167.98 ng	5094770	1,4-Dichlorobenzene-d4	6.73

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		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
5		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9



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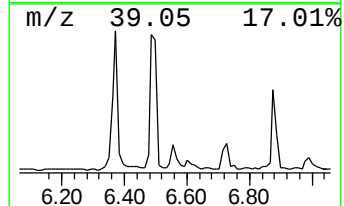
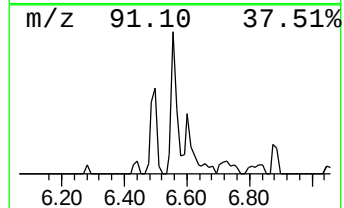
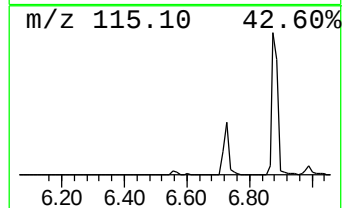
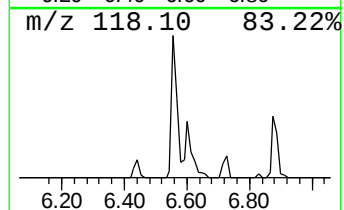
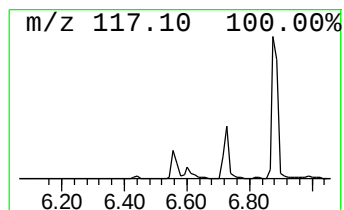
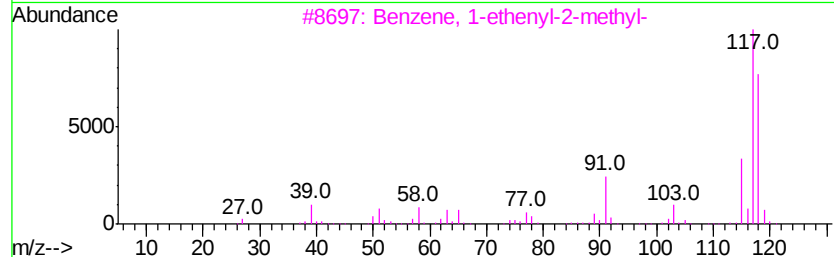
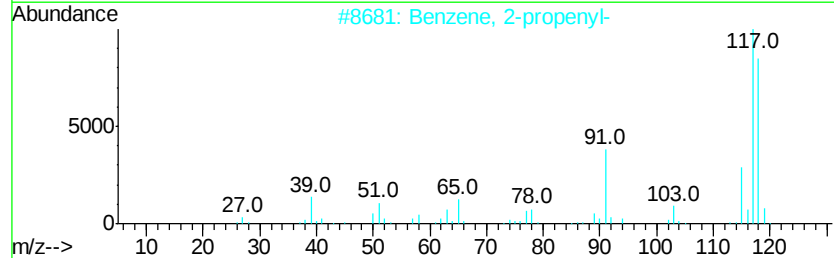
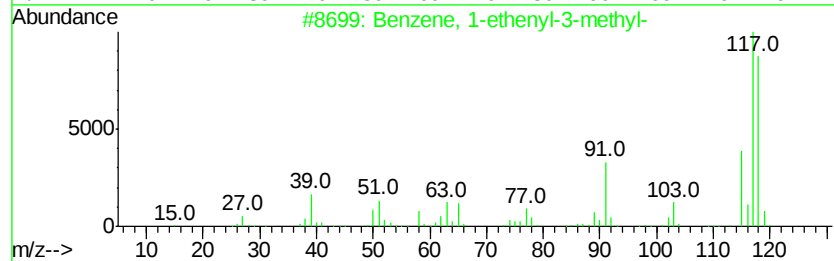
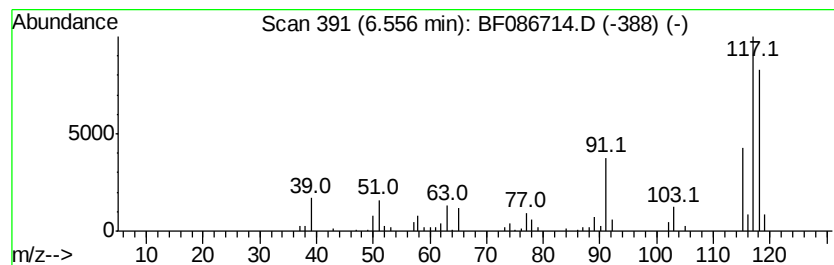
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethenyl-3-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.56	5.07 ng	153729	1,4-Dichlorobenzene-d4	6.73	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	96
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	94
3	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	93
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	93
5	Bicyclo[4.2.0]octa-1,3,5-triene, ...	118	C9H10	022250-74-4	92



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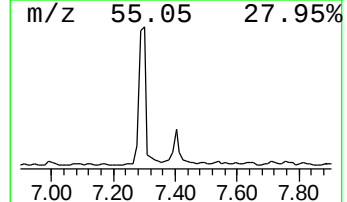
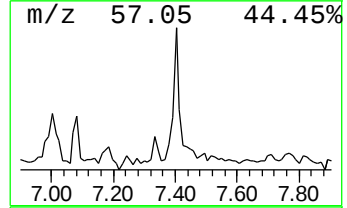
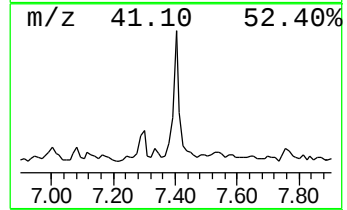
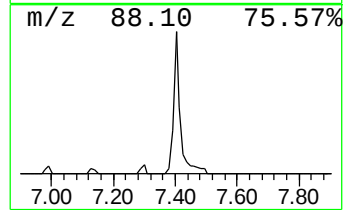
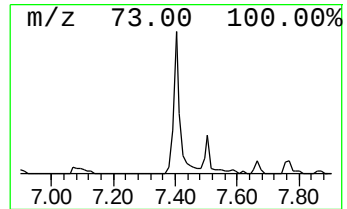
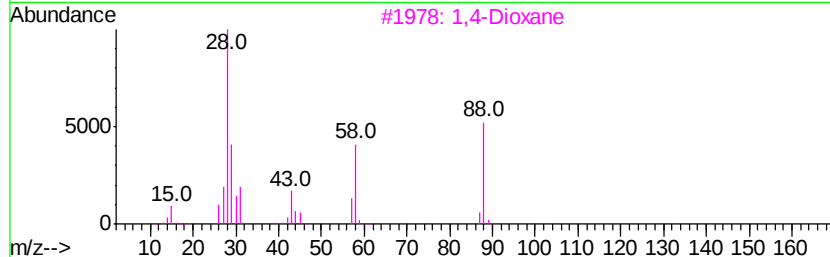
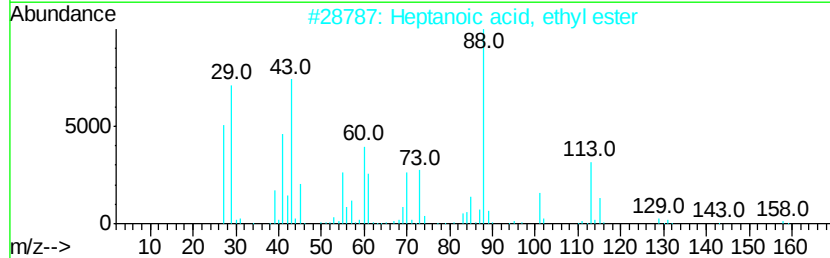
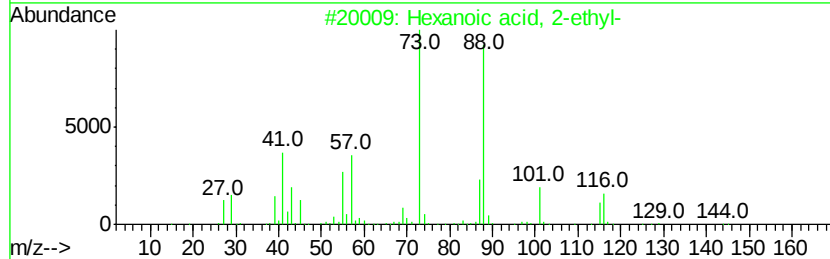
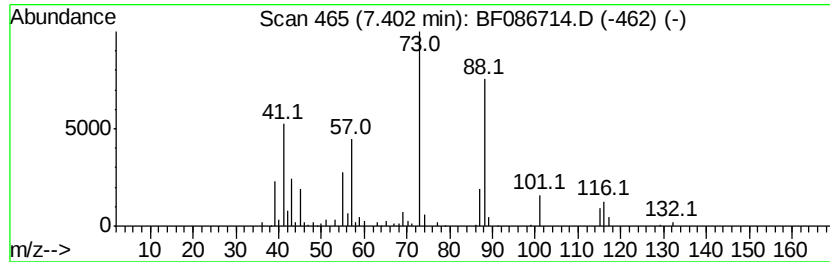
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Peak Number 5 Hexanoic acid, 2-ethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.40	4.59 ng	176578	Naphthalene-d8	8.01
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexanoic acid, 2-ethyl-	144 C8H16O2	000149-57-5	78
2	Heptanoic acid, ethyl ester	158 C9H18O2	000106-30-9	39
3	1,4-Dioxane	88 C4H8O2	000123-91-1	28
4	Cyclohexaneacetic acid, .alpha.-...	170 C10H18O2	006051-00-9	28
5	Pentanoic acid, 2-methyl-, methy...	130 C7H14O2	002177-77-7	25



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EUT01-TD-B12(0-8.5)-C

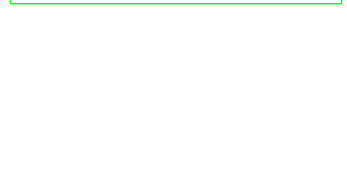
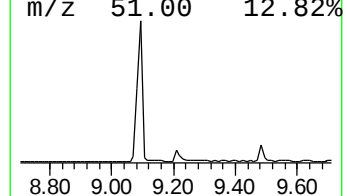
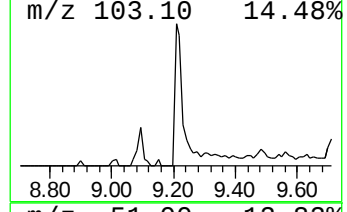
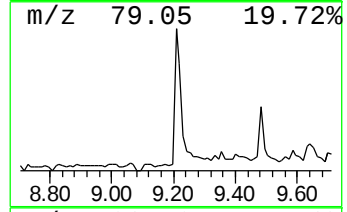
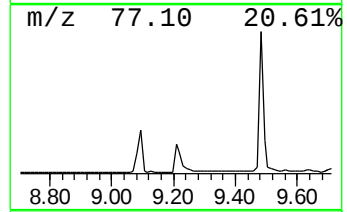
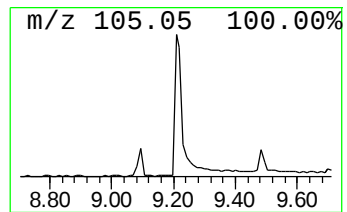
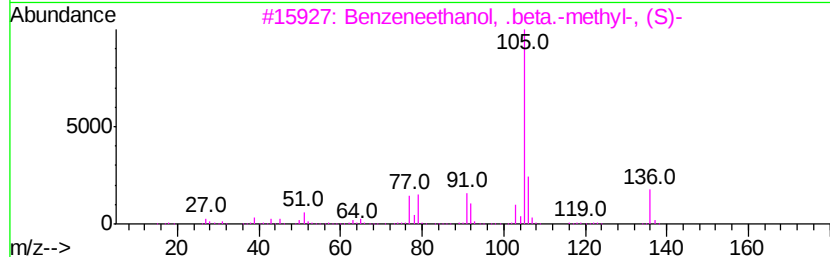
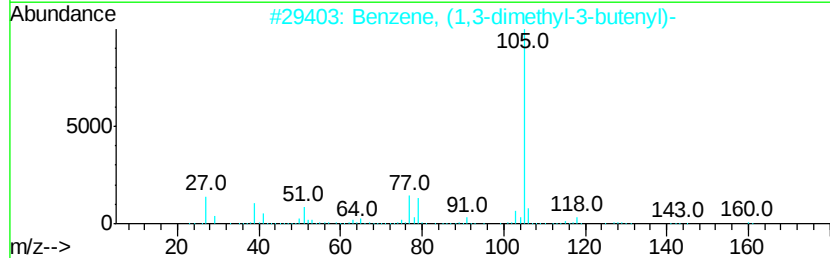
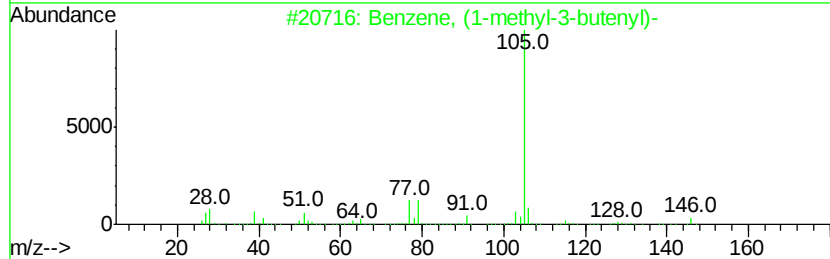
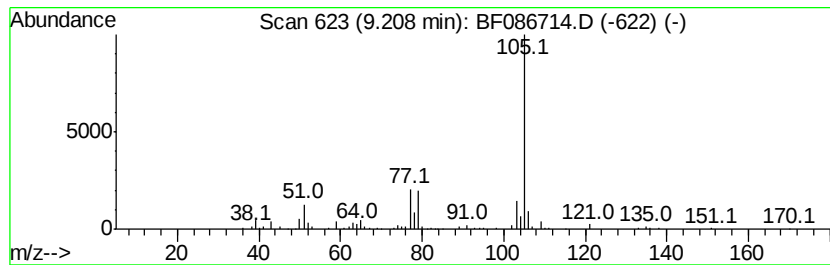
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, (1-methyl-3-butenyl)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.21	5.07 ng	209403	Acenaphthene-d10	9.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-3-butenyl)-	146	C11H14	010340-49-5	72
2		Benzene, (1,3-dimethyl-3-butenyl)-	160	C12H16	056851-51-5	72
3		Benzeneethanol, .beta.-methyl-, ...	136	C9H12O	037778-99-7	64
4		Benzenemethanethiol, 4-methyl-	138	C8H10S	004498-99-1	64
5		Benzene, (1-nitroethyl)-	151	C8H9NO2	007214-61-1	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050516\
Data File : BF086714.D
Acq On : 5 May 2016 19:21
Operator : UM/SJ
Sample : H2829-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EUT01-TD-B12(0-8.5)-C

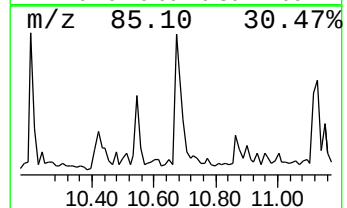
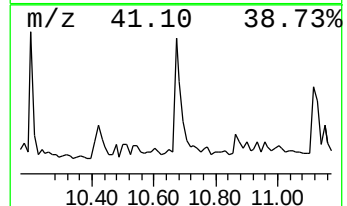
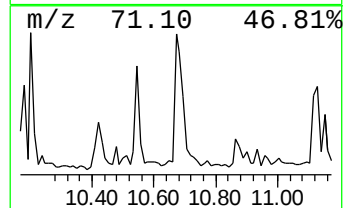
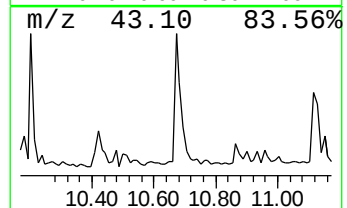
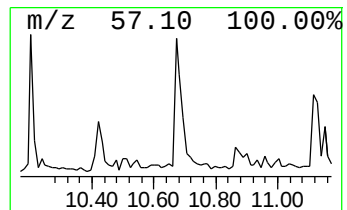
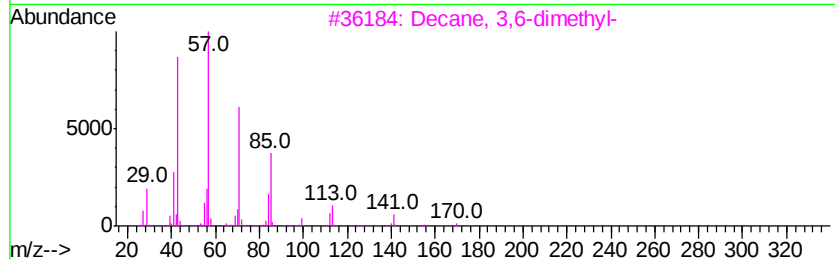
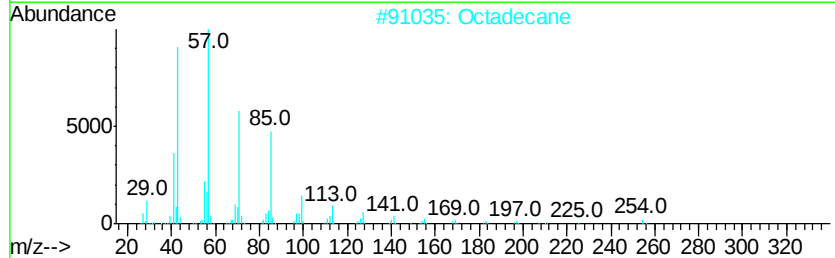
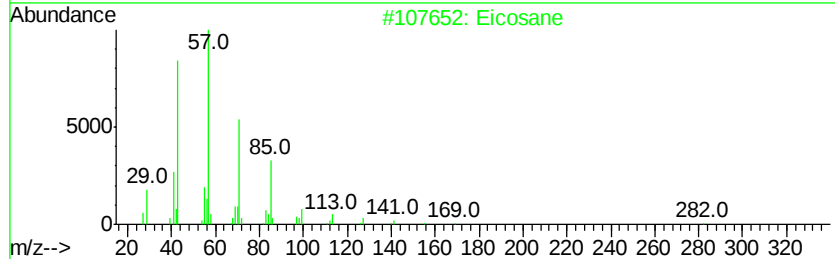
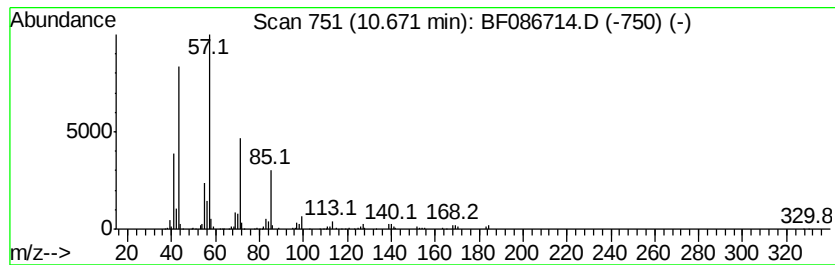
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.67	7.37 ng	279538	Phenanthrene-d10	11.24

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	91
2	Octadecane		254	C18H38	000593-45-3	90
3	Decane, 3,6-dimethyl-		170	C12H26	017312-53-7	87
4	Hexadecane		226	C16H34	000544-76-3	87
5	1-Iodo-2-methylundecane		296	C12H25I	073105-67-6	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050516\
Data File : BF086714.D
Acq On : 5 May 2016 19:21
Operator : UM/SJ
Sample : H2829-01
Misc :
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Instrument :
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ClientSampleId :
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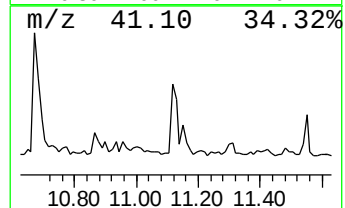
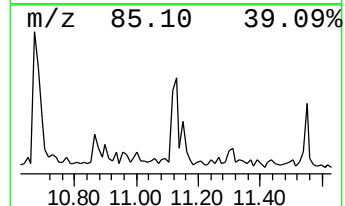
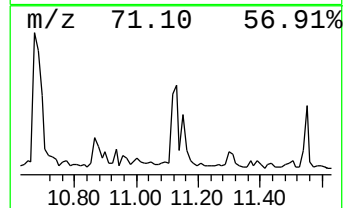
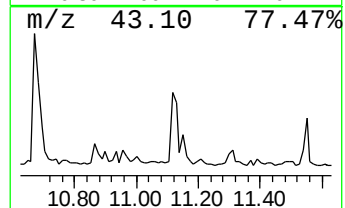
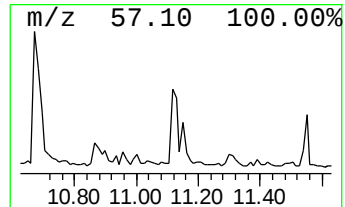
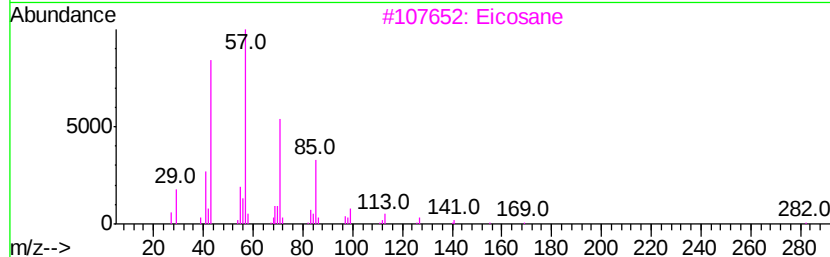
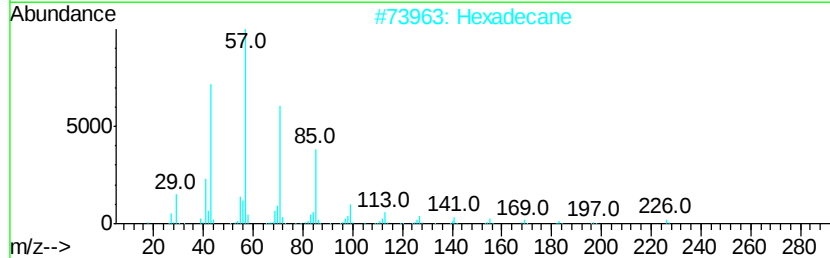
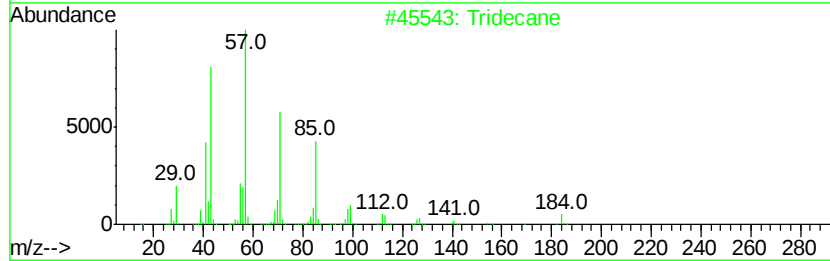
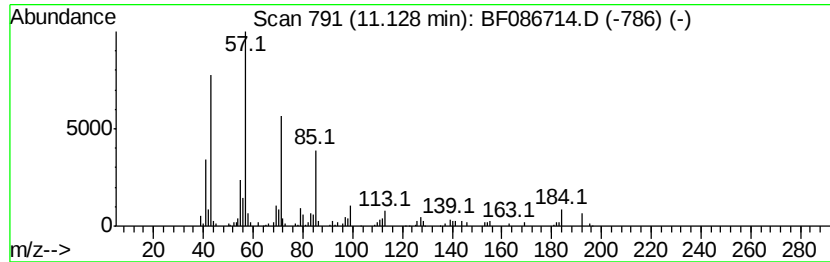
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Tridecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.13	7.28 ng	276253	Phenanthrene-d10	11.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	90
2		Hexadecane	226	C16H34	000544-76-3	87
3		Eicosane	282	C20H42	000112-95-8	87
4		Dodecane, 3-methyl-	184	C13H28	017312-57-1	87
5		Heptadecane	240	C17H36	000629-78-7	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050516\
Data File : BF086714.D
Acq On : 5 May 2016 19:21
Operator : UM/SJ
Sample : H2829-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EUT01-TD-B12(0-8.5)-C

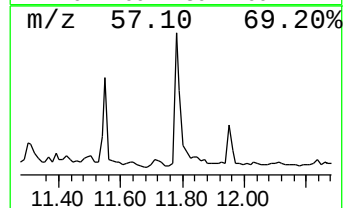
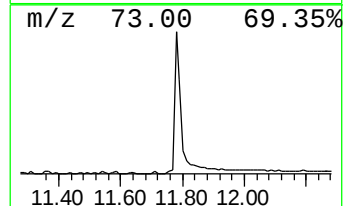
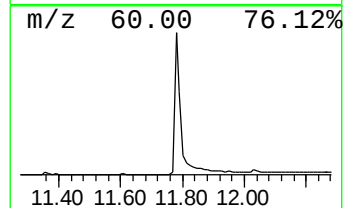
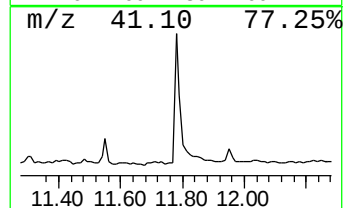
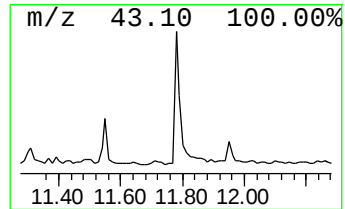
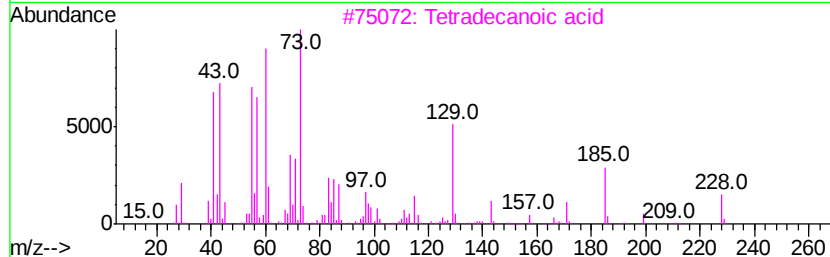
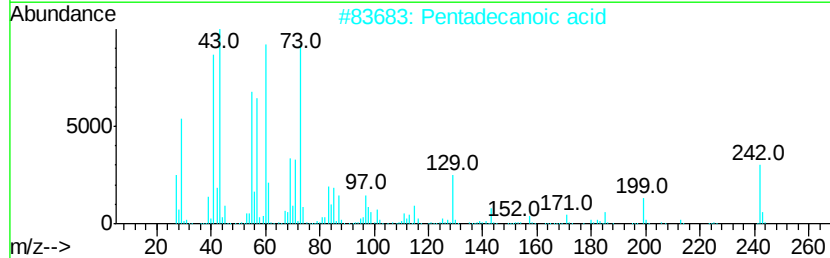
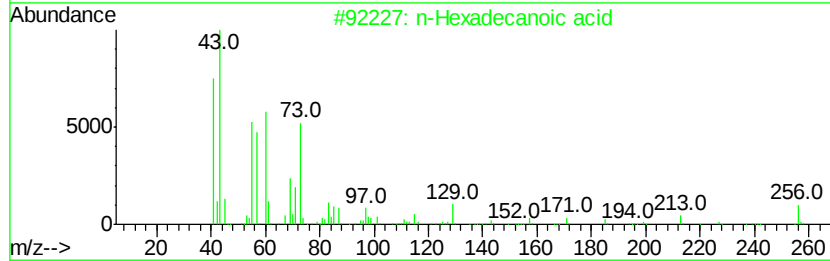
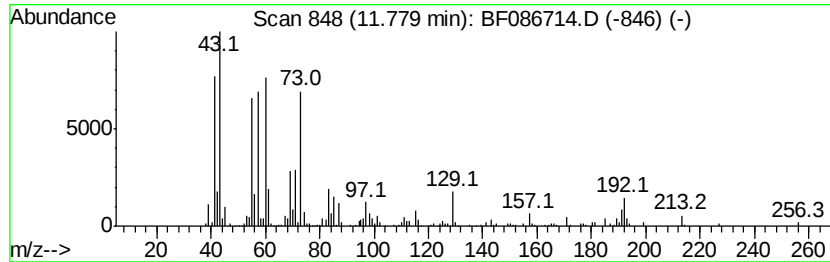
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.78	12.49 ng	473789	Phenanthrene-d10	11.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4		Undecanoic acid	186	C11H22O2	000112-37-8	64
5		Tridecanoic acid	214	C13H26O2	000638-53-9	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050516\
Data File : BF086714.D
Acq On : 5 May 2016 19:21
Operator : UM/SJ
Sample : H2829-01
Misc :
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Instrument :
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ClientSampleId :
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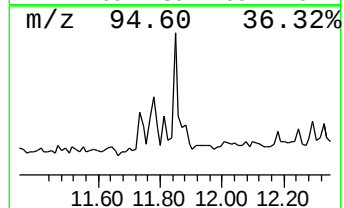
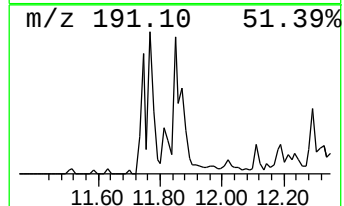
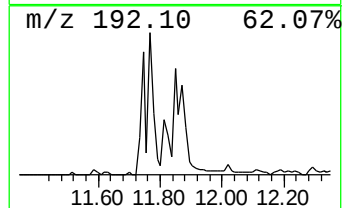
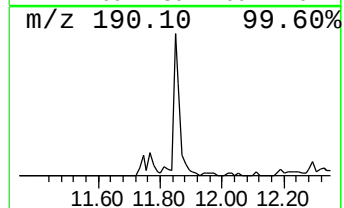
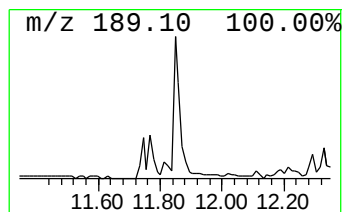
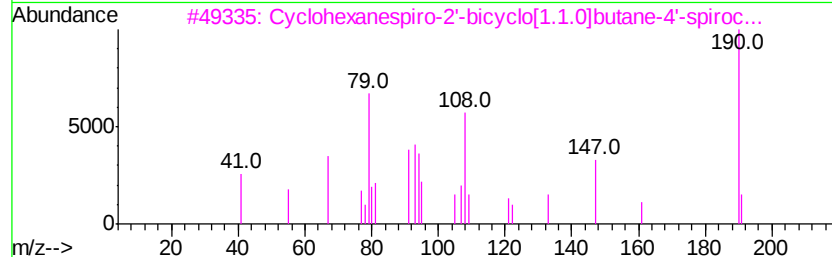
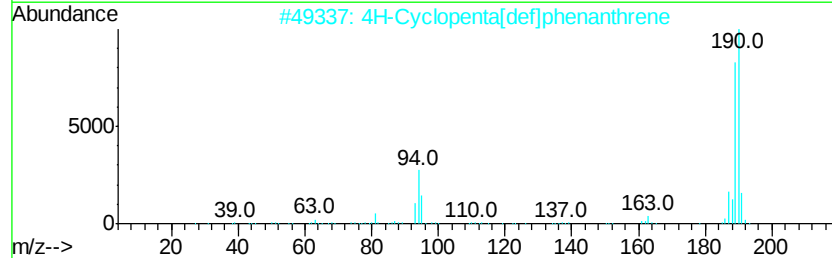
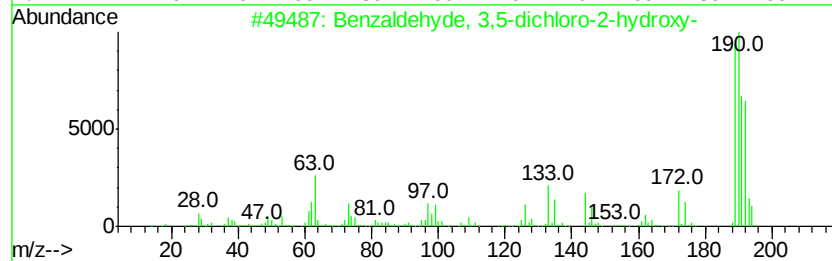
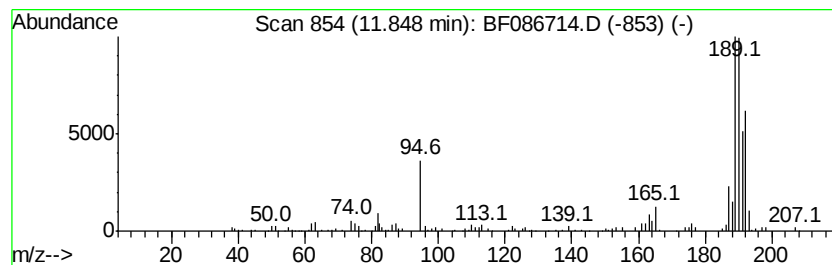
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzaldehyde, 3,5-dichloro-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.85	5.13 ng	194753	Phenanthrene-d10	11.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehydve, 3,5-dichloro-2-hvd...	190	C7H4Cl2O2	000090-60-8	72
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
3			Cyclohexanespiro-2'-bicyclo[1.1.0]butane-4'-spiro...	190	C14H22	107924-94-7	22
4			Benzene, 1,1'-(2-cyclopropen-1-yl)-	192	C15H12	022825-21-4	22
5			Benzene, 1,1'-(1,2-propadienyldi-	192	C15H12	014251-57-1	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050516\
Data File : BF086714.D
Acq On : 5 May 2016 19:21
Operator : UM/SJ
Sample : H2829-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

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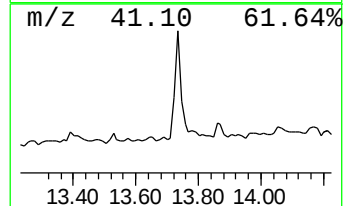
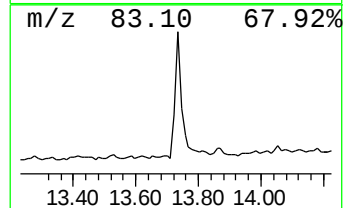
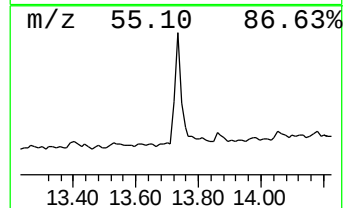
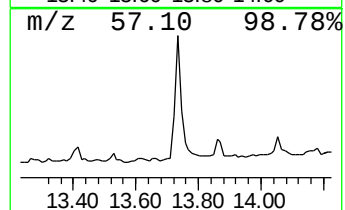
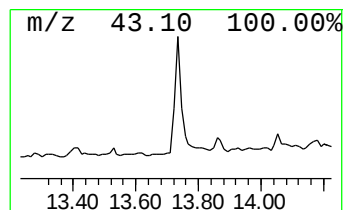
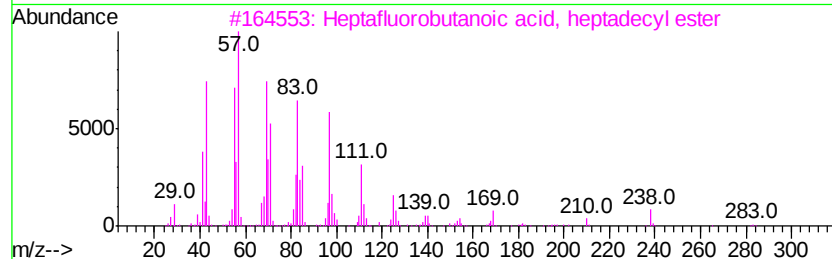
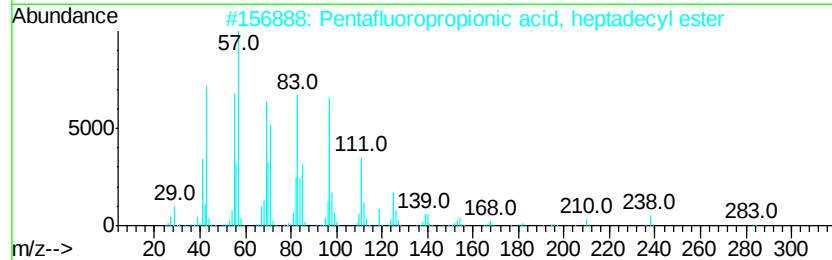
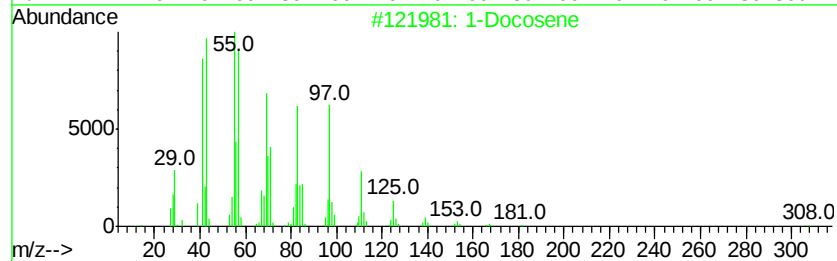
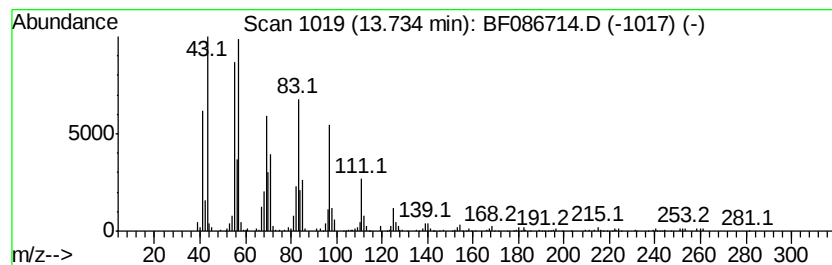
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 1-Docosene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	7.13 ng	299085	Chrysene-d12	13.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	95
2		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	94
3		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	91
4		Trifluoroacetic acid, n-octadecy...	366	C20H37F3O2	079392-43-1	90
5		Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050516\
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Acq On : 5 May 2016 19:21
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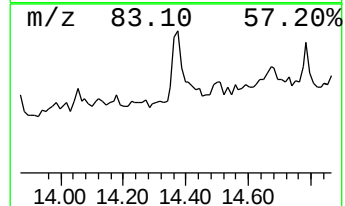
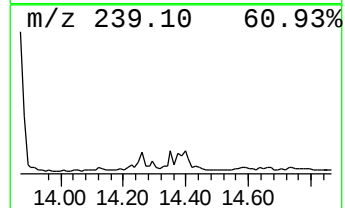
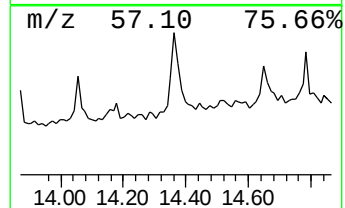
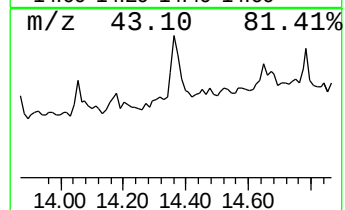
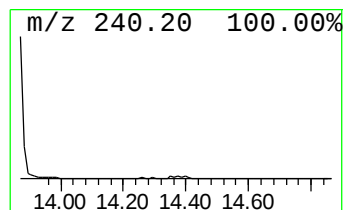
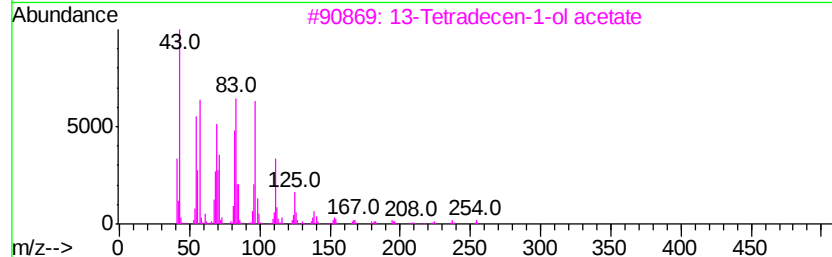
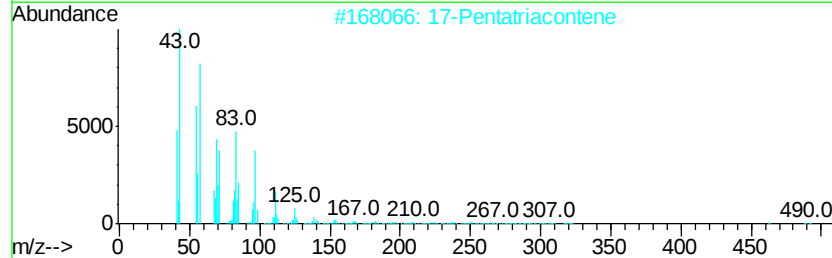
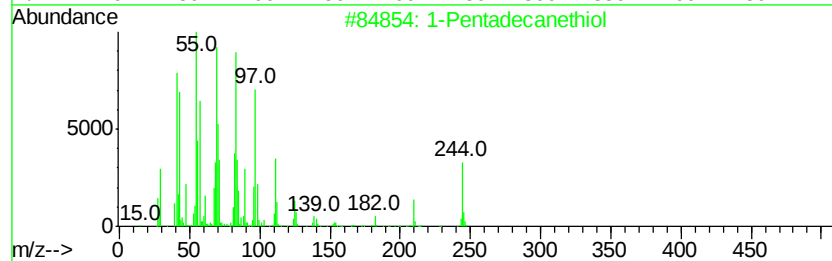
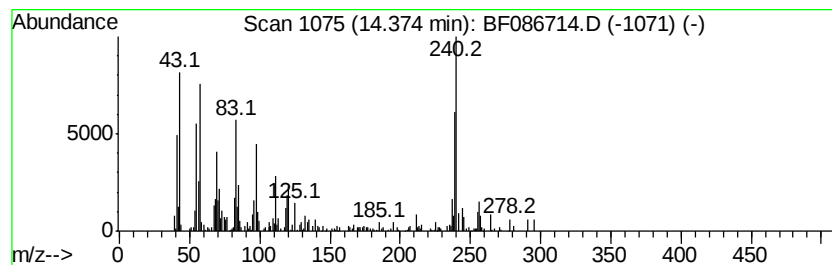
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown14.37 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.37	4.09 ng	171516	Chrysene-d12	13.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentadecanethiol	244	C15H32S	025276-70-4	27
2			17-Pentatriacontene	491	C35H70	006971-40-0	27
3			13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	25
4			2-Methyl-E-7-hexadecene	238	C17H34	064183-52-4	25
5			Undecane, 1-cyclohexyl-3-(2-cycl...	348	C25H48	007225-69-6	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050516\
Data File : BF086714.D
Acq On : 5 May 2016 19:21
Operator : UM/SJ
Sample : H2829-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EUT01-TD-B12(0-8.5)-C

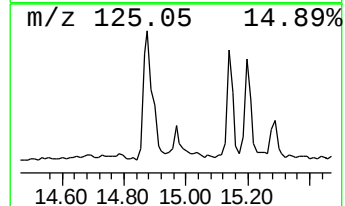
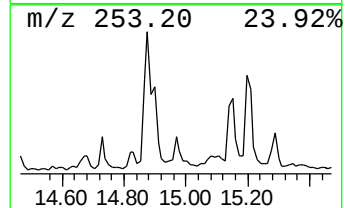
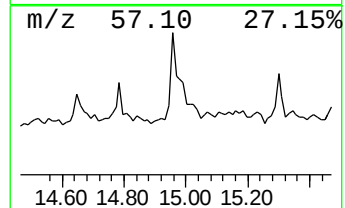
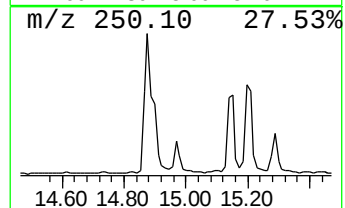
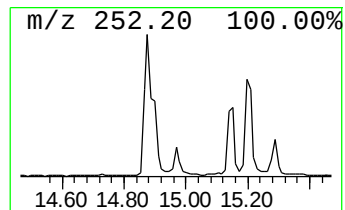
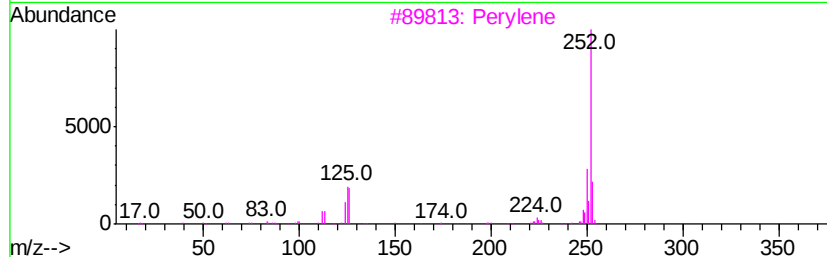
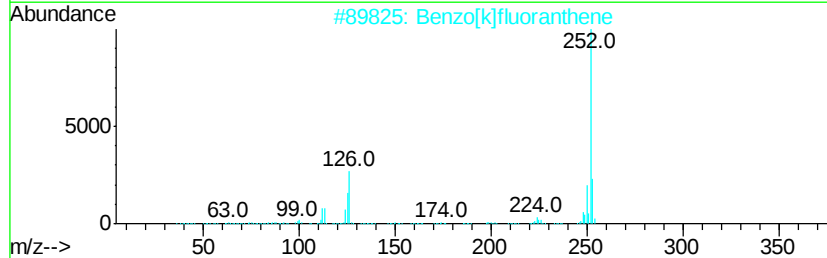
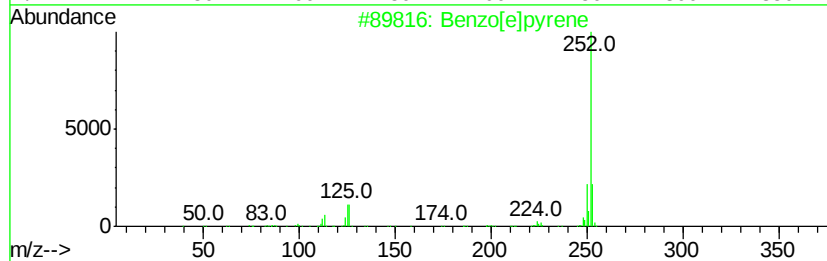
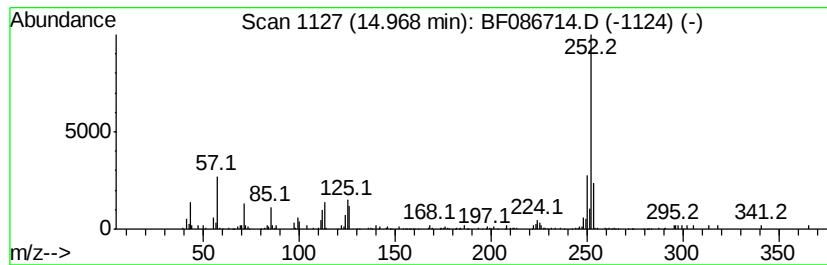
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.97	7.75 ng	181624	Perylene-d12	15.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	91
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	90
3			Perylene	252	C20H12	000198-55-0	90
4			Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	90
5			Benzo[j]fluoranthene	252	C20H12	000205-82-3	89



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050516\
Data File : BF086714.D
Acq On : 5 May 2016 19:21
Operator : UM/SJ
Sample : H2829-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

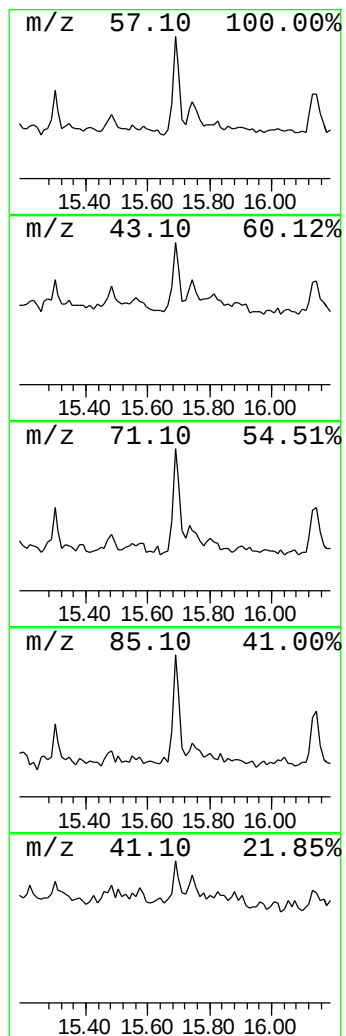
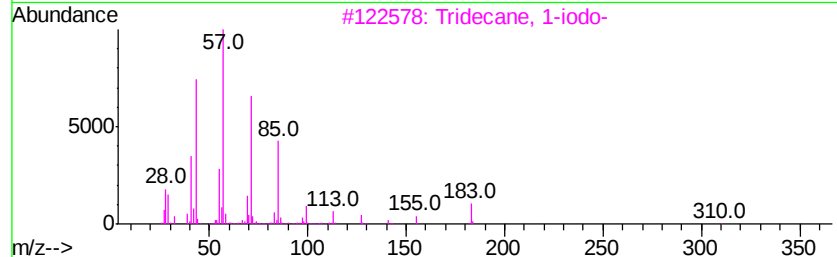
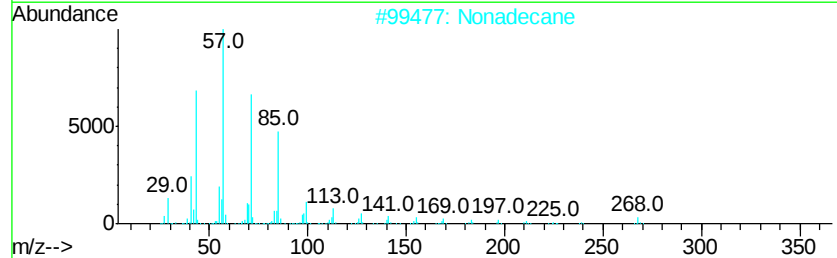
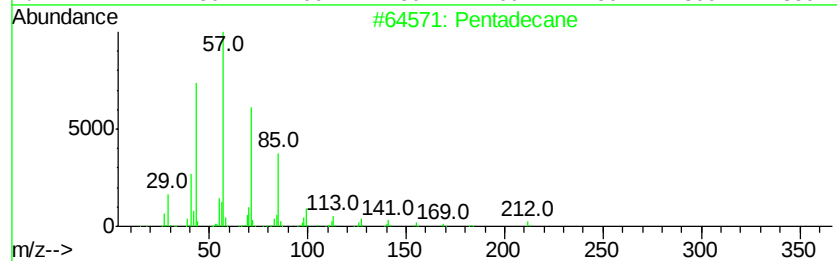
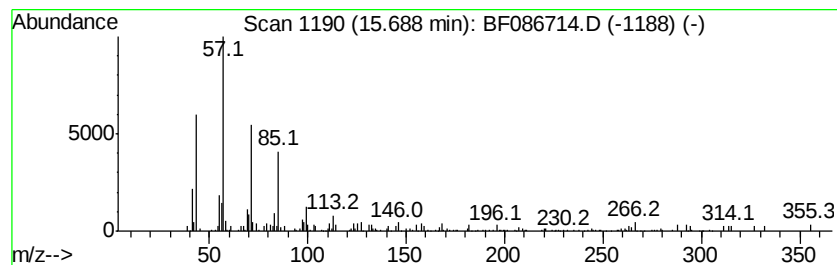
Instrument :
BNA_F
ClientSampleId :
EUT01-TD-B12(0-8.5)-C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Pentadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.69	4.16 ng	97481	Perylene-d12		15.27
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane	212	C15H32	000629-62-9	80
2	Nonadecane	268	C19H40	000629-92-5	80
3	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	80
4	Tetracosane	338	C24H50	000646-31-1	80
5	Tridecane, 6-propyl-	226	C16H34	055045-10-8	76



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050516\
 Data File : BF086714.D
 Acq On : 5 May 2016 19:21
 Operator : UM/SJ
 Sample : H2829-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EUT01-TD-B12(0-8.5)-C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.89	16.1 ng		489078	1	6.73	1213180	40.0
unknown6.50	6.50	168.0 ng		5094770	1	6.73	1213180	40.0
Benzene, 1-etheny...	6.56	5.1 ng		153729	1	6.73	1213180	40.0
Hexanoic acid, 2-...	7.40	4.6 ng		176578	2	8.01	1540000	40.0
Benzene, (1-methy...	9.21	5.1 ng		209403	3	9.76	1650980	40.0
Eicosane	10.67	7.4 ng		279538	4	11.24	1517900	40.0
Tridecane	11.13	7.3 ng		276253	4	11.24	1517900	40.0
n-Hexadecanoic acid	11.78	12.5 ng		473789	4	11.24	1517900	40.0
Benzaldehyde, 3,5...	11.85	5.1 ng		194753	4	11.24	1517900	40.0
1-Docosene	13.73	7.1 ng		299085	5	13.87	1677250	40.0
unknown14.37	14.37	4.1 ng		171516	5	13.87	1677250	40.0
Benzo[e]pyrene	14.97	7.8 ng		181624	6	15.27	937424	40.0
Pentadecane	15.69	4.2 ng		97481	6	15.27	937424	40.0