

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061416\
 Data File : BF088020.D
 Acq On : 15 Jun 2016 1:50
 Operator : UM/SJ
 Sample : H3471-07
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 STA-1037-(0-4)

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-BF060716.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	21	rVB	225466	372576	13.63%	1.193%
2	1.850	21	23	34	rVV	1131857	1844242	67.48%	5.905%
3	1.998	34	36	69	rVB3	96774	389084	14.24%	1.246%
4	4.959	293	295	302	rBV	126971	190579	6.97%	0.610%
5	5.382	329	332	334	rBV	2217446	2281573	83.49%	7.305%
6	6.422	420	423	426	rBV	1698510	2495243	91.30%	7.990%
7	6.547	432	434	437	rBV	1828789	2501657	91.54%	8.010%
8	6.616	438	440	442	rBV	126506	101955	3.73%	0.326%
9	6.776	452	454	457	rBV	406549	559864	20.49%	1.793%
10	6.936	466	468	471	rVB	1855255	1886678	69.04%	6.041%
11	7.347	501	504	507	rBV	1347626	1510637	55.28%	4.837%
12	7.462	510	514	520	rBV	31811	49500	1.81%	0.158%
13	8.068	565	567	571	rBV	855677	749574	27.43%	2.400%
14	9.153	659	662	664	rBV	2684566	2732905	100.00%	8.751%
15	9.542	694	696	698	rBV	487445	476209	17.43%	1.525%
16	9.679	707	708	711	rBV	46616	53314	1.95%	0.171%
17	9.759	714	715	718	rBV	44008	48012	1.76%	0.154%
18	9.828	718	721	722	rBV	750208	777296	28.44%	2.489%
19	10.262	758	759	761	rVB	93551	72763	2.66%	0.233%
20	10.479	775	778	781	rBV	32867	56986	2.09%	0.182%
21	10.616	787	790	792	rBV	1204246	1412452	51.68%	4.523%
22	10.731	799	800	807	rBV	77164	121388	4.44%	0.389%
23	11.188	835	840	845	rVB3	34054	85998	3.15%	0.275%
24	11.302	848	850	854	rBV2	697504	890771	32.59%	2.852%
25	11.382	854	857	858	rBV	93974	98133	3.59%	0.314%
26	11.531	868	870	875	rBV5	22512	52155	1.91%	0.167%
27	11.805	892	894	895	rBV	37892	33606	1.23%	0.108%
28	11.839	895	897	899	rVB	37202	48073	1.76%	0.154%
29	11.919	902	904	908	rVB	74296	111536	4.08%	0.357%
30	12.011	908	912	917	rBV5	23329	59222	2.17%	0.190%
31	12.102	917	920	921	rBV	37803	38443	1.41%	0.123%
32	12.354	939	942	944	rBV	64622	76142	2.79%	0.244%
33	12.388	944	945	948	rVV3	31364	59092	2.16%	0.189%
34	12.434	948	949	954	rVB2	39699	55719	2.04%	0.178%

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35	12.514	954	956	959	rBV	722256	630350	23.07%	2.018%
36	12.662	966	969	973	rBV4	21899	54048	1.98%	0.173%
37	12.742	973	976	978	rVV	700471	772912	28.28%	2.475%
38	12.799	978	981	983	rVB3	47405	75232	2.75%	0.241%
39	12.857	984	986	987	rBV	27746	34496	1.26%	0.110%
40	12.891	987	989	992	rVB	1870665	2023923	74.06%	6.480%
41	12.959	992	995	999	rBV3	68036	115832	4.24%	0.371%
42	13.074	1001	1005	1006	rVB2	85301	161531	5.91%	0.517%
43	13.142	1006	1011	1012	rBV2	45844	85366	3.12%	0.273%
44	13.177	1012	1014	1015	rBV	93053	106132	3.88%	0.340%
45	13.268	1020	1022	1023	rVB	55787	48550	1.78%	0.155%
46	13.302	1023	1025	1027	rBV2	32212	51773	1.89%	0.166%
47	13.428	1034	1036	1037	rBV	64663	67664	2.48%	0.217%
48	13.474	1037	1040	1043	rVB5	37650	78631	2.88%	0.252%
49	13.577	1045	1049	1051	rVV	84658	112983	4.13%	0.362%
50	13.691	1056	1059	1060	rVV2	61712	104018	3.81%	0.333%
51	13.714	1060	1061	1062	rVV	68171	60461	2.21%	0.194%
52	13.737	1062	1063	1066	rVV2	58895	86256	3.16%	0.276%
53	13.794	1066	1068	1072	rVB3	84204	156976	5.74%	0.503%
54	13.942	1077	1081	1087	rBV2	744895	1505879	55.10%	4.822%
55	14.045	1087	1090	1091	rBV2	42624	59530	2.18%	0.191%
56	14.114	1094	1096	1098	rVB3	26127	37401	1.37%	0.120%
57	14.320	1112	1114	1116	rVB	57105	73819	2.70%	0.236%
58	14.354	1116	1117	1119	rVB2	44133	53996	1.98%	0.173%
59	14.422	1119	1123	1124	rBV3	50115	101891	3.73%	0.326%
60	14.525	1129	1132	1134	rVB3	22473	35612	1.30%	0.114%
61	14.628	1139	1141	1142	rBV	27494	36746	1.34%	0.118%
62	14.720	1146	1149	1152	rBV3	32398	70268	2.57%	0.225%
63	14.788	1152	1155	1159	rBV2	38275	91644	3.35%	0.293%
64	14.948	1166	1169	1174	rBV	292803	560544	20.51%	1.795%
65	15.051	1174	1178	1180	rVV3	59141	102825	3.76%	0.329%
66	15.234	1191	1194	1197	rVB	137308	191132	6.99%	0.612%
67	15.291	1197	1199	1202	rVV	176223	219006	8.01%	0.701%
68	15.348	1202	1204	1211	rVB	456285	624285	22.84%	1.999%
69	15.794	1241	1243	1245	rVB2	21129	32735	1.20%	0.105%
70	16.388	1293	1295	1302	rVB6	19789	49134	1.80%	0.157%
71	16.525	1304	1307	1309	rBV2	13041	35586	1.30%	0.114%

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Integrator: RTE
Smoothing : ON Filtering: 5
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72	16.697	1319	1322	1328	rVB2	66298	126724	4.64%	0.406%
73	17.108	1354	1358	1363	rVB	46113	102062	3.73%	0.327%

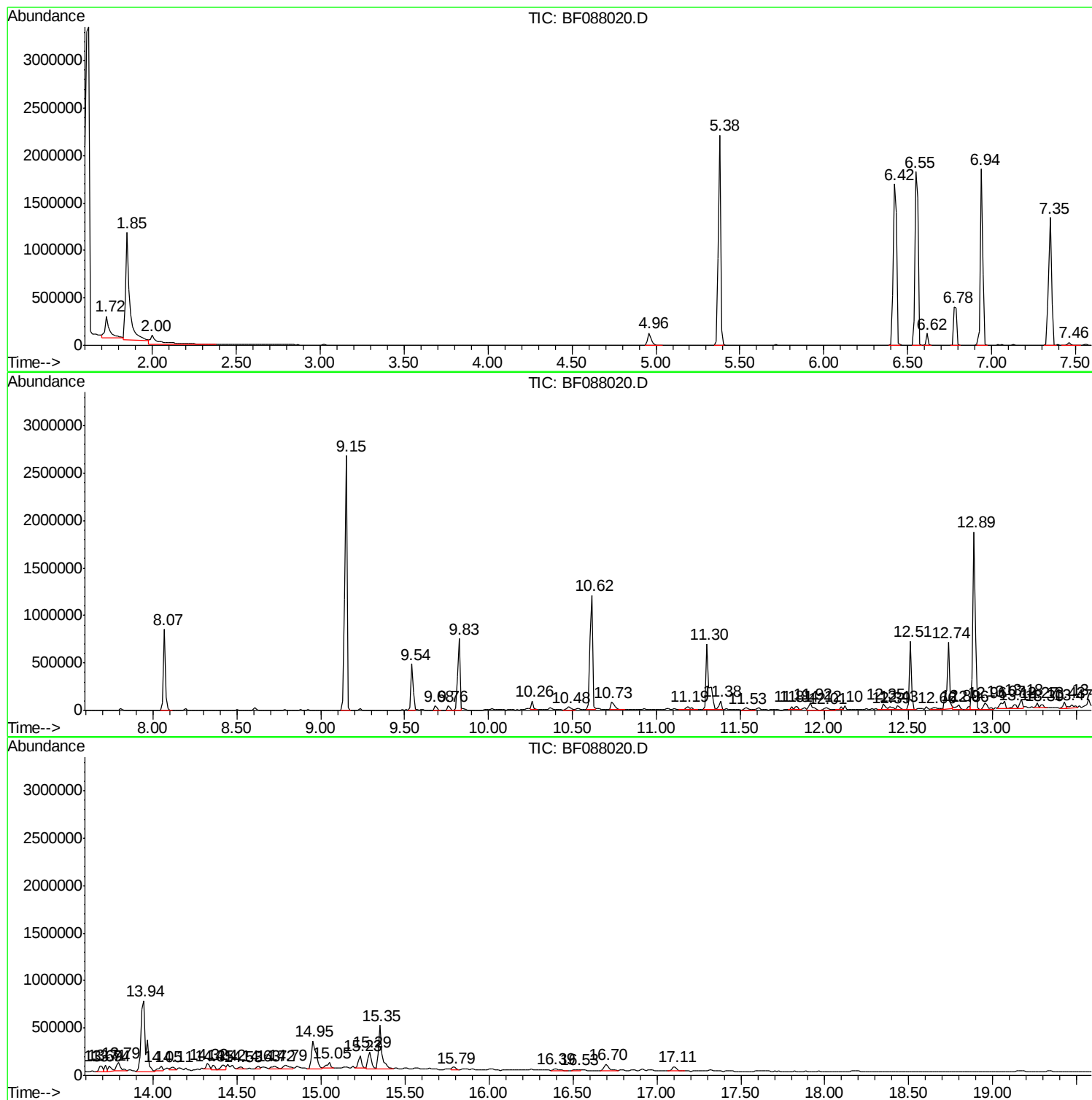
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Instrument :
BNA_F
Client Sampled :
STA-1037-(0-4)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.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 : H3471-07
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Instrument :
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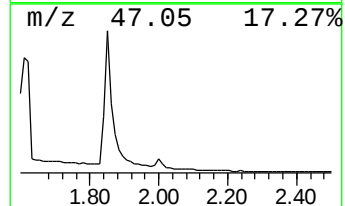
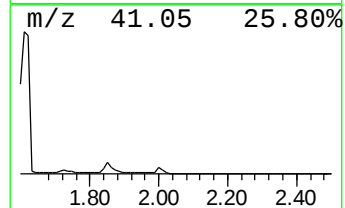
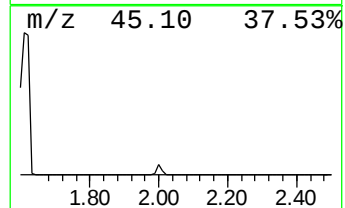
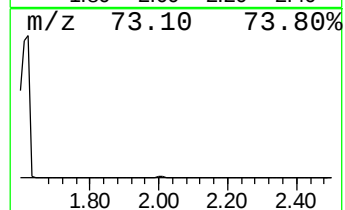
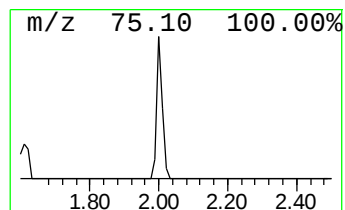
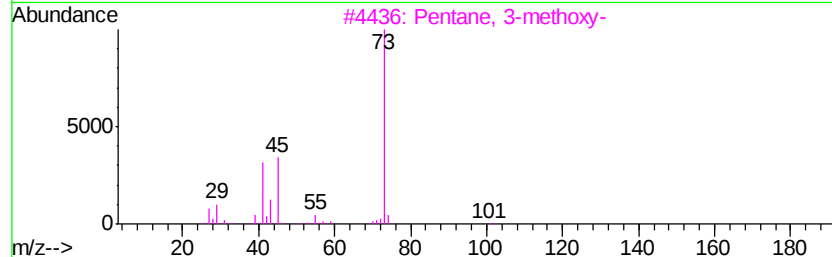
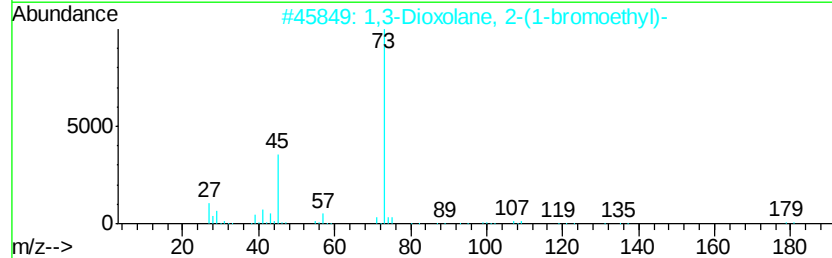
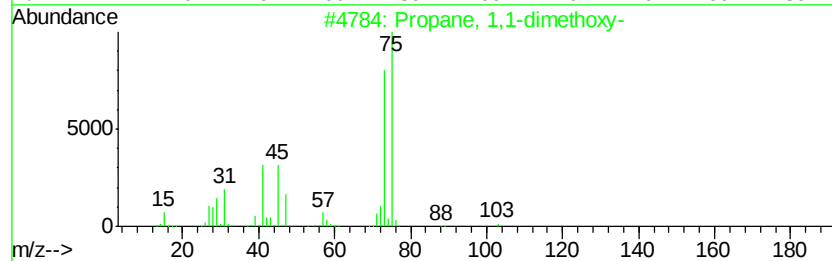
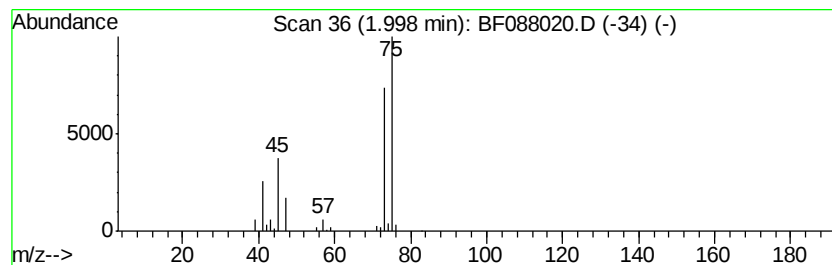
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Propane, 1,1-dimethoxy- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.00	13.90 ng	389084	1,4-Dichlorobenzene-d4	6.79

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	83	
2	1,3-Dioxolane, 2-(1-bromoethyl)-	180	C5H9BrO2	005267-73-2	12	
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12	
4	Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9	
5	Glycine	75	C2H5NO2	000056-40-6	9	



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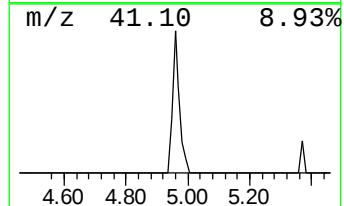
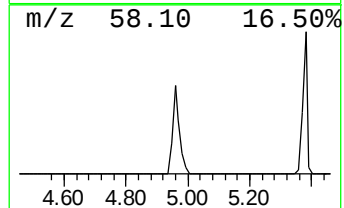
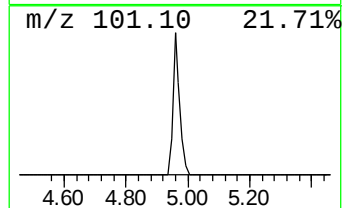
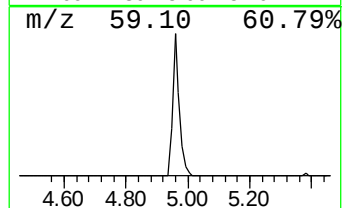
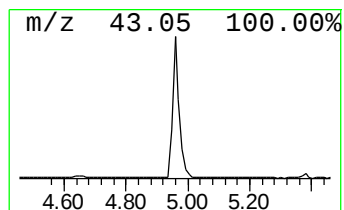
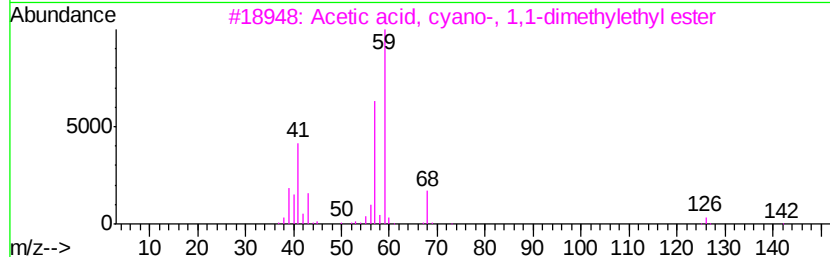
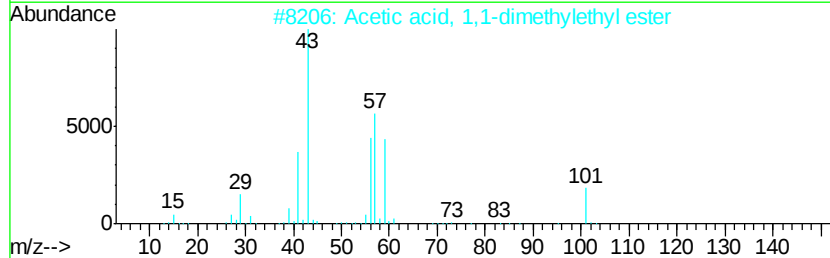
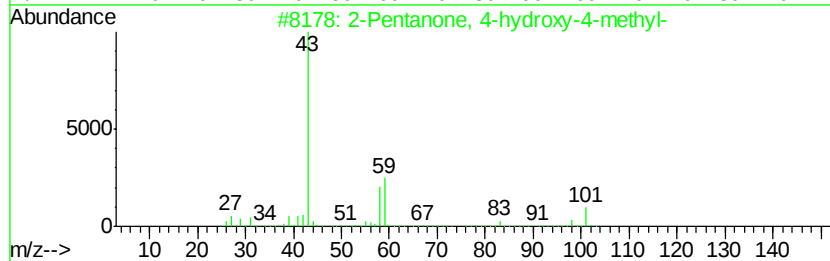
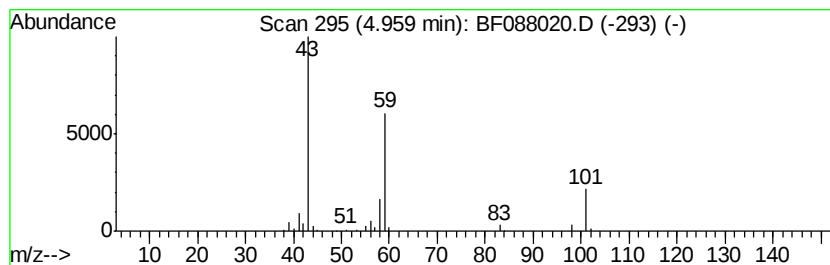
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	6.81 ng	190579	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	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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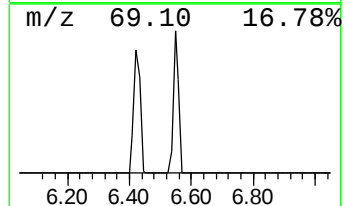
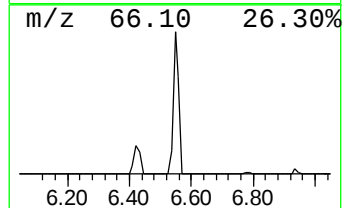
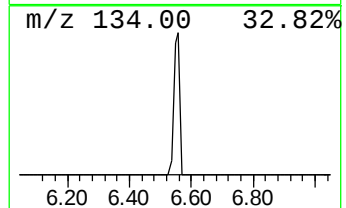
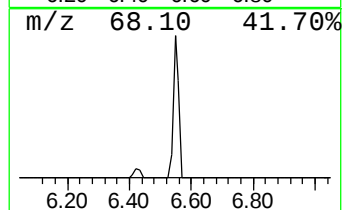
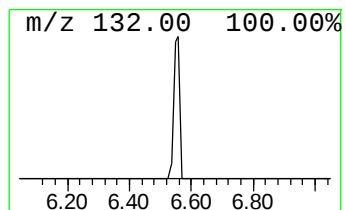
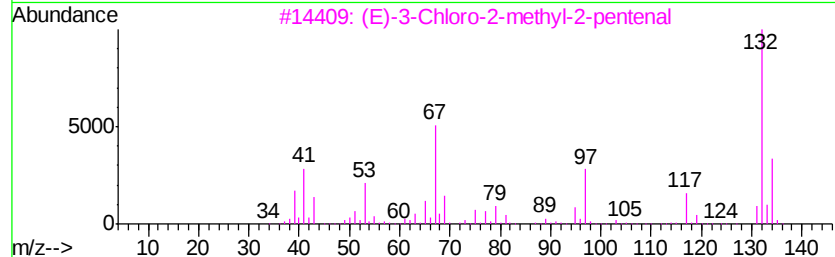
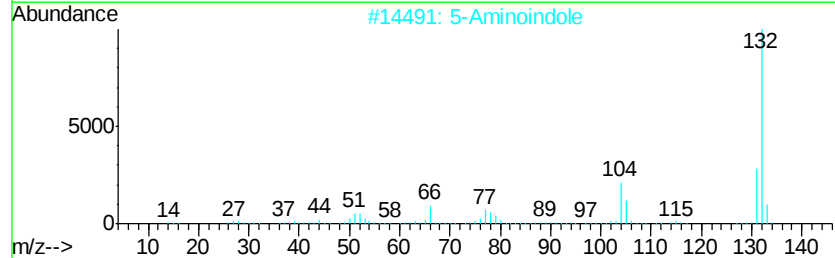
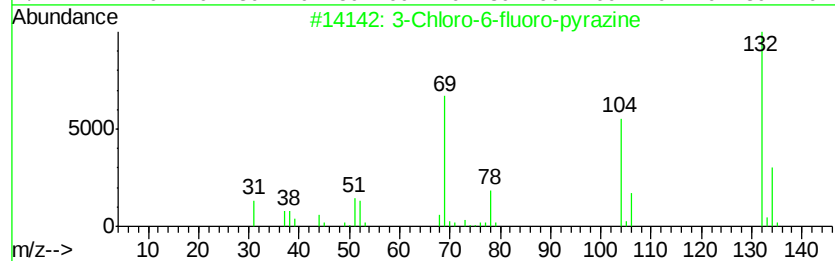
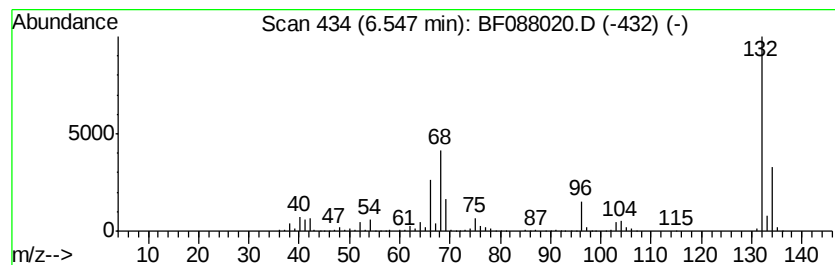
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TIC Library : C:\Database\NIST11.L
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Peak Number 5 unknown6.55 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.55	89.37 ng	2501660	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27	
2	5-Aminoindole	132	C8H8N2	005192-03-0	12	
3	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9	
4	1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9	
5	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9	



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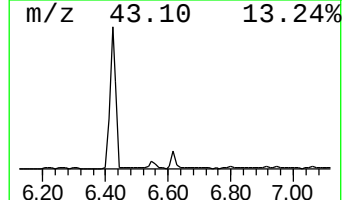
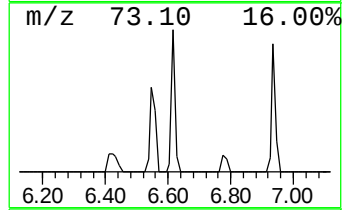
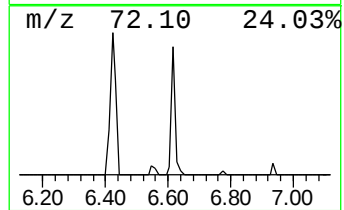
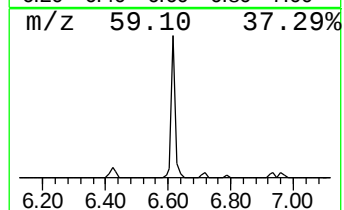
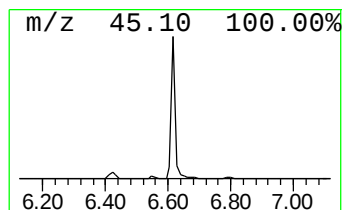
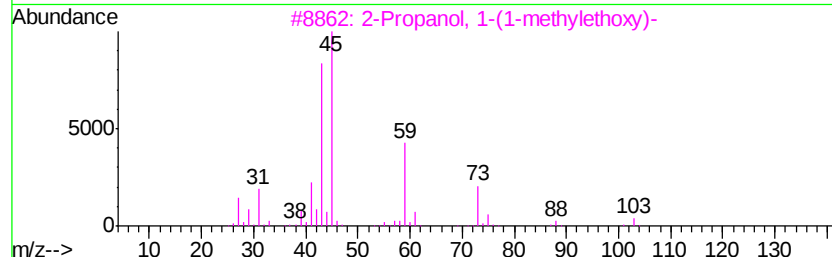
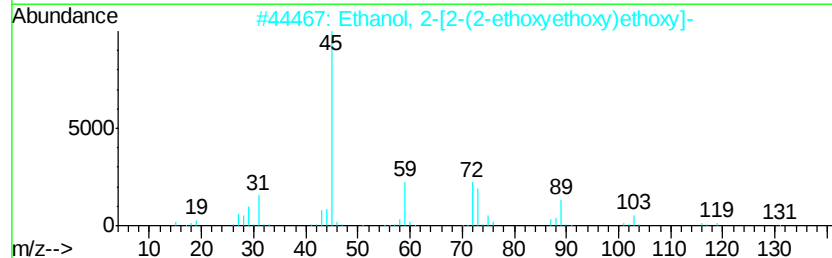
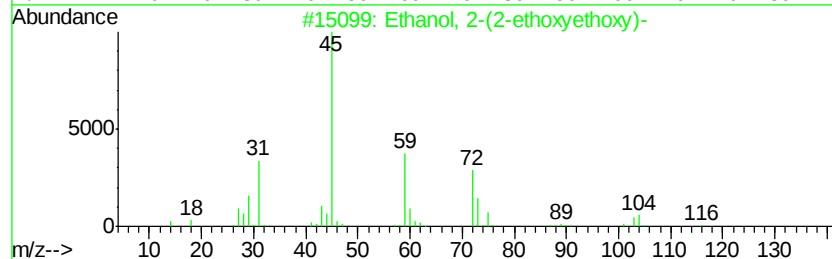
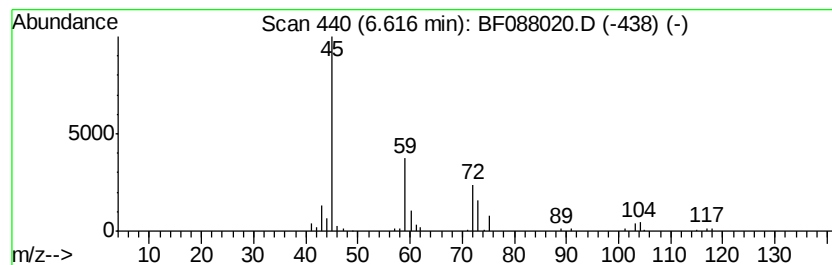
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	3.64 ng	101955	1,4-Dichlorobenzene-d4	6.79

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ethanol, 2-(2-ethoxvethoxy)-	134	C6H14O3	000111-90-0	90
2		Ethanol, 2-[2-(2-ethoxvethoxy)et...	178	C8H18O4	000112-50-5	64
3		2-Propanol, 1-(1-methvlethoxy)-	118	C6H14O2	003944-36-3	45
4		Ethanol, 2-(2-methoxvethoxy)-	120	C5H12O3	000111-77-3	42
5		Diethyl carbitol	162	C8H18O3	000112-36-7	34



Data Path : Z:\HPCHEM1\BNA F\DATA\BF061416\
Data File : BF088020.D
Acq On : 15 Jun 2016 1:50
Operator : UM/SJ
Sample : H3471-07
Misc :
ALS Vial : 24 Sample Multiplier: 1

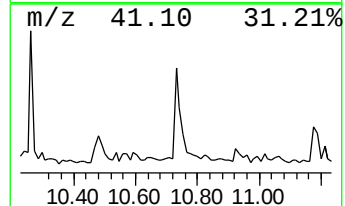
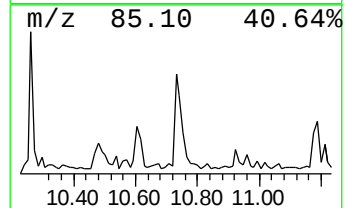
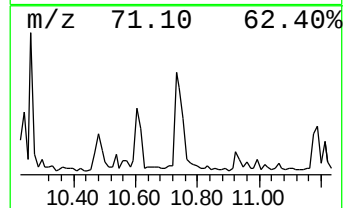
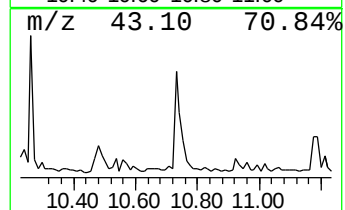
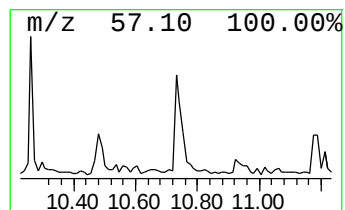
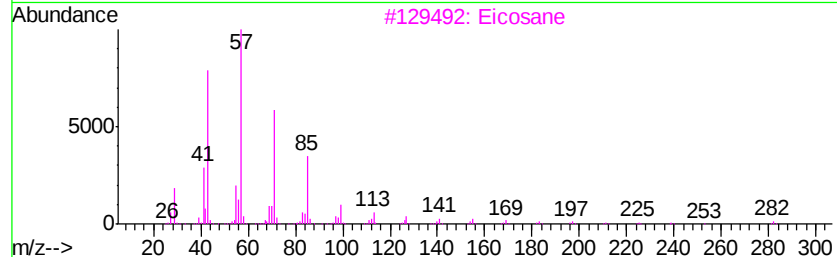
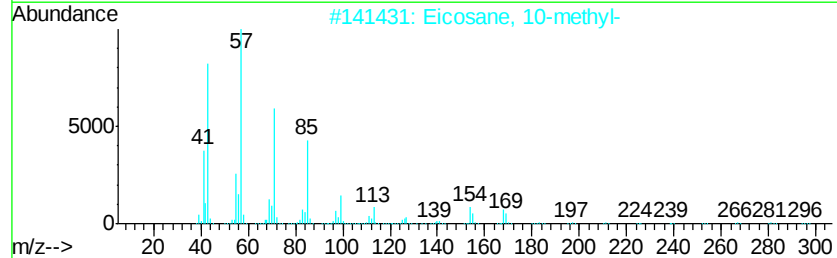
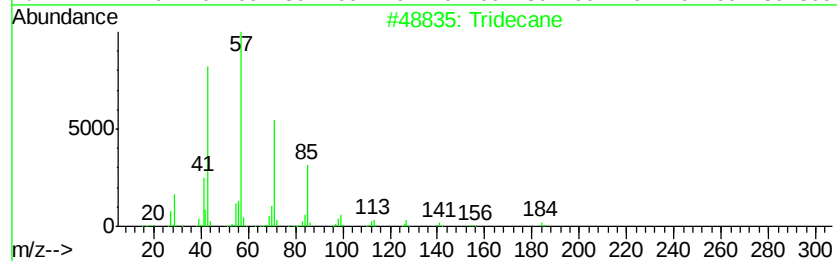
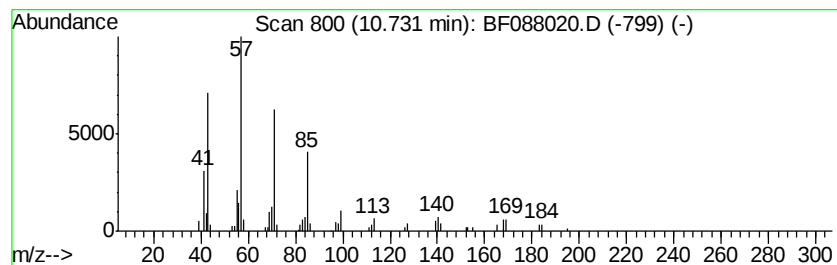
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Tridecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.73	2.73 ng	121388	Phenanthrene-d10		11.30	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane	184	C13H28	000629-50-5	95
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	83
3		Eicosane	282	C20H42	000112-95-8	80
4		Heptadecane	240	C17H36	000629-78-7	80
5		Hexadecane	226	C16H34	000544-76-3	80



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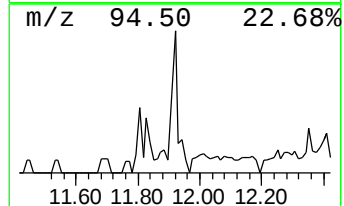
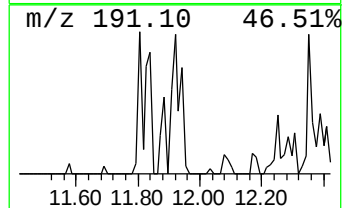
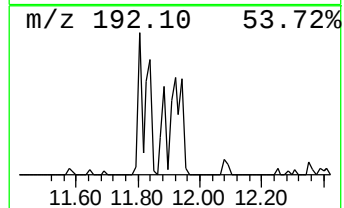
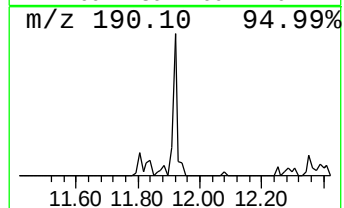
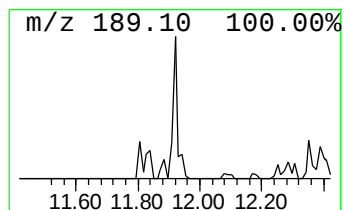
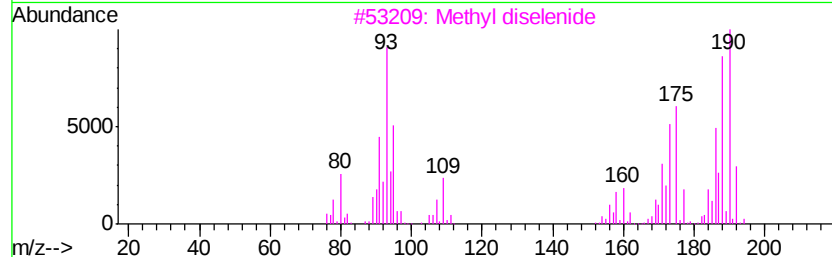
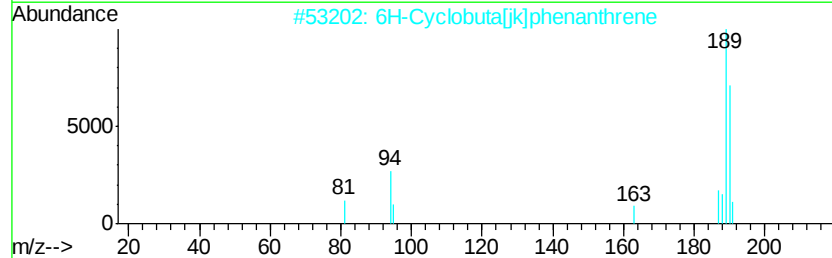
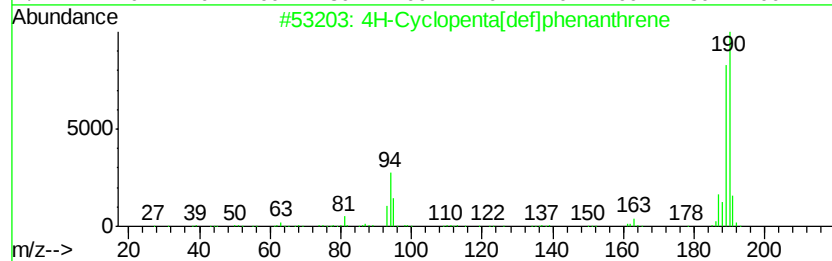
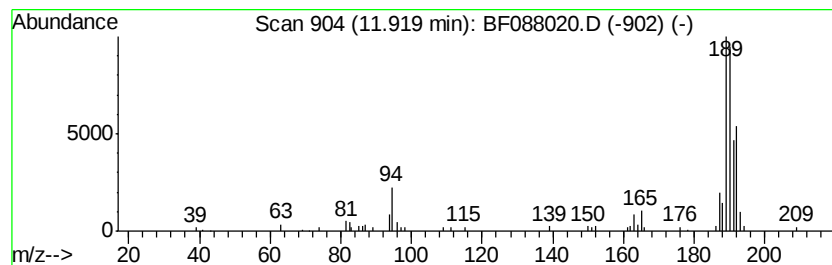
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.92	2.50 ng	111536	Phenanthrene-d10	11.30	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2	6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	58
3	Methyl diselenide	190	C2H6Se2	007101-31-7	41
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	35
5	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	30



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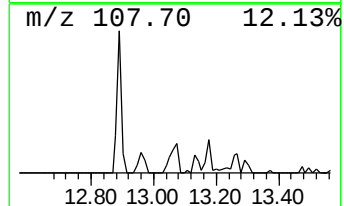
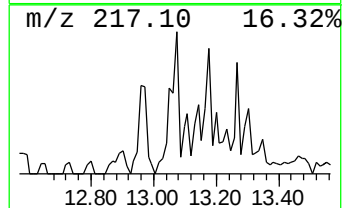
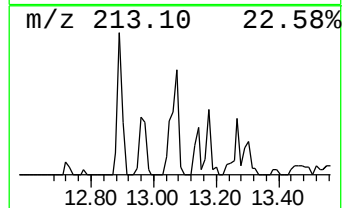
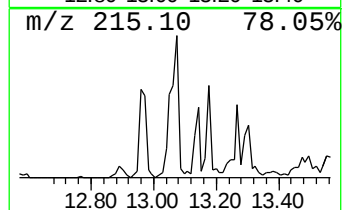
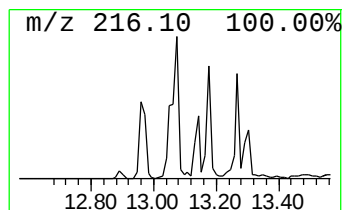
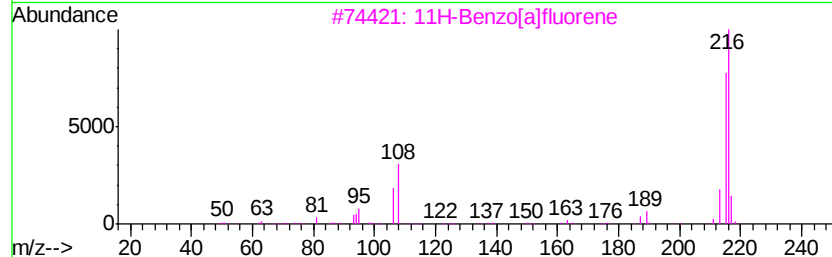
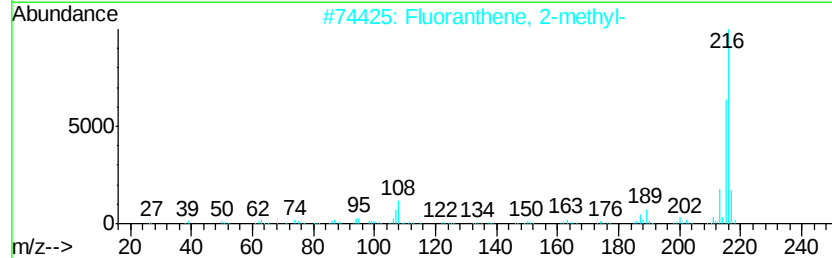
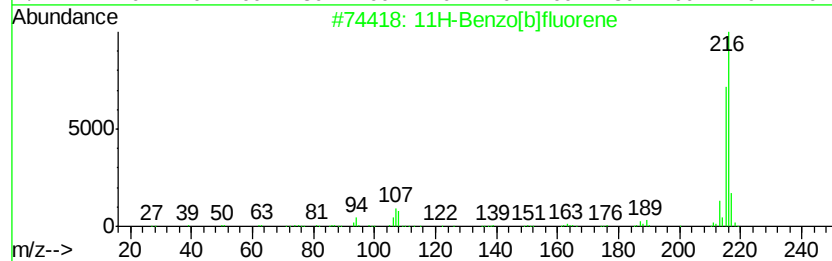
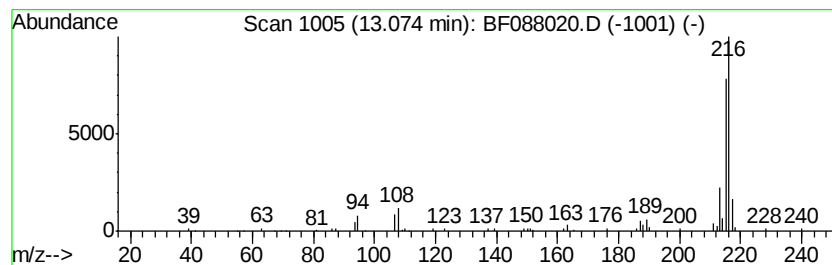
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.07	2.15 ng	161531	Chrysene-d12	13.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	95
2			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	94
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
4			Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
5			Pyrene, 2-methyl-	216	C17H12	003442-78-2	87



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Data File : BF088020.D
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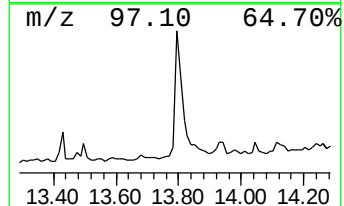
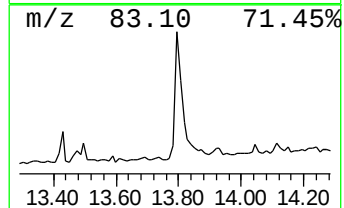
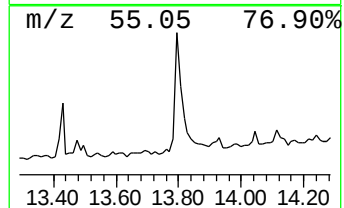
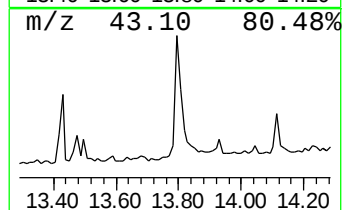
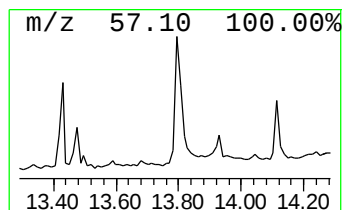
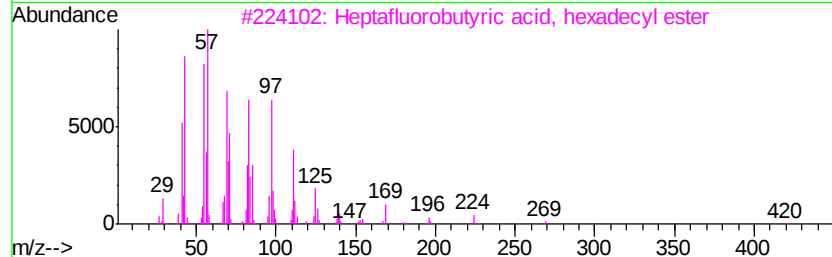
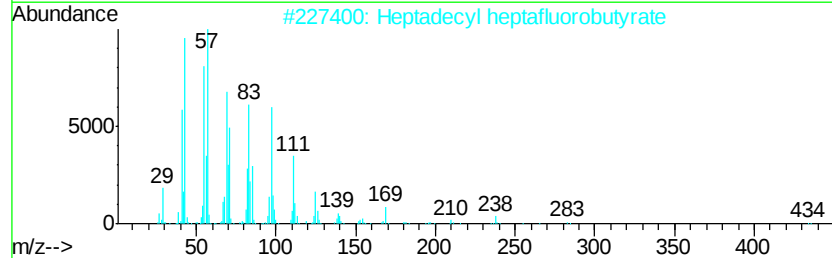
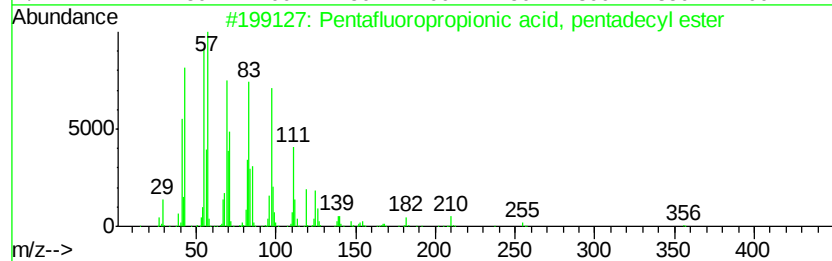
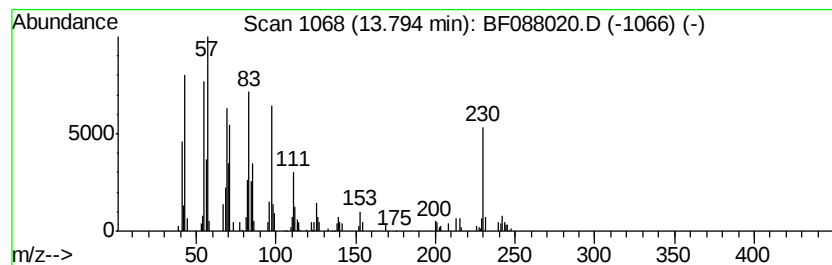
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Pentafluoropropionic acid, ... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	2.08 ng	156976	Chrysene-d12	13.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	92
2		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	89
3		Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	89
4		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	86
5		Octacosyl heptafluorobutyrate	606	C32H57F7O2	1000351-83-6	76



Data Path : Z:\HPCHEM1\BNA F\DATA\BF061416\
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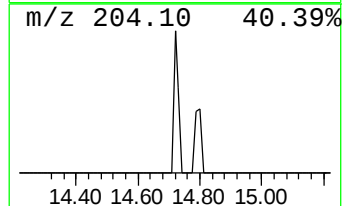
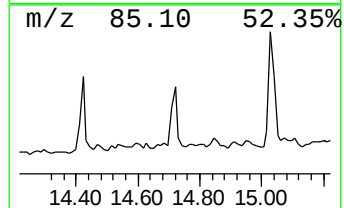
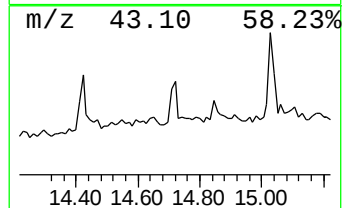
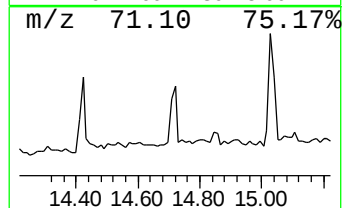
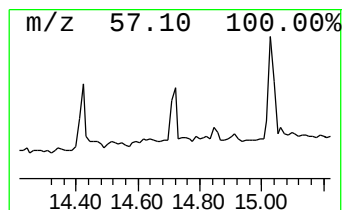
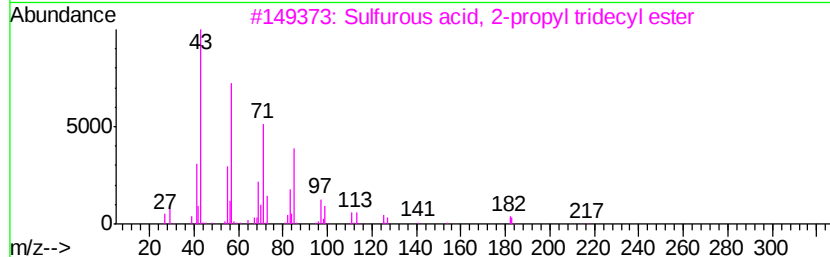
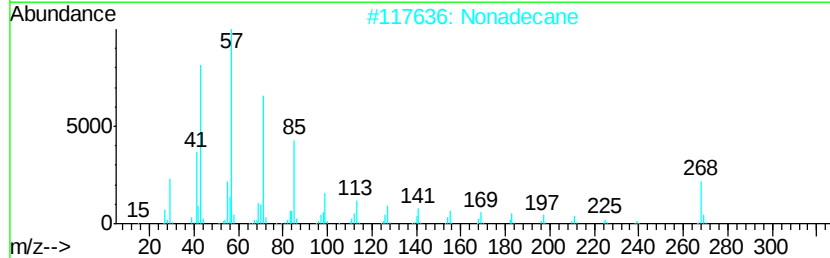
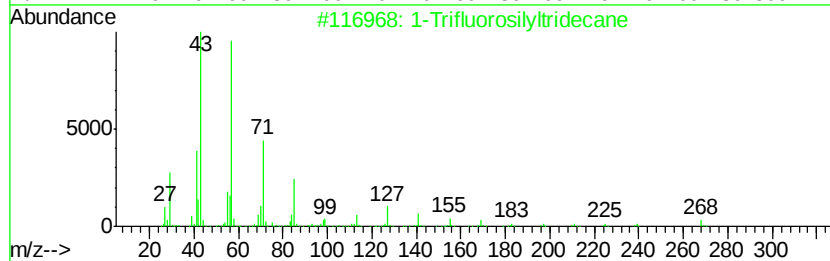
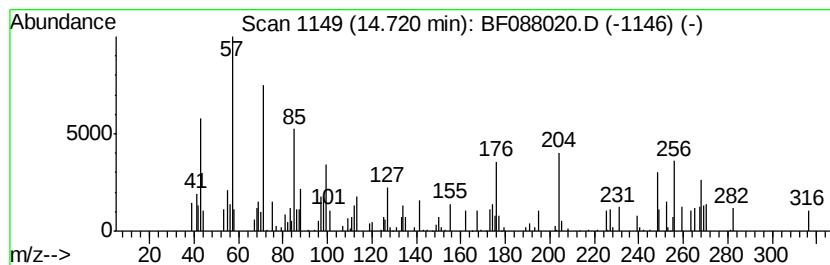
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Trifluorosilyltridecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.72	2.25 ng	70268	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Trifluorosilyltridecane	268	C13H27F3Si	1000217-08-2	50
2		Nonadecane	268	C19H40	000629-92-5	45
3		Sulfurous acid, 2-propyl tridecyl...	306	C16H34O3S	1000309-12-4	35
4		Decane, 2,3,8-trimethyl-	184	C13H28	062238-14-6	35
5		Pentadecane, 2-methyl-	226	C16H34	001560-93-6	35



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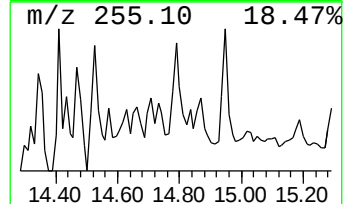
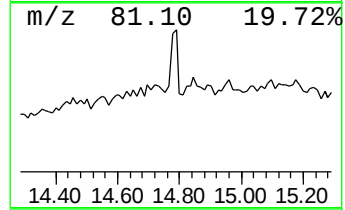
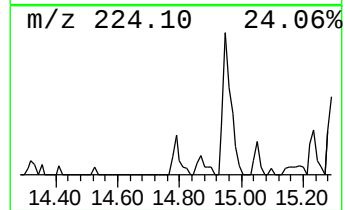
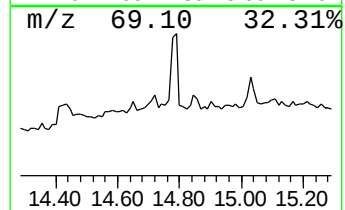
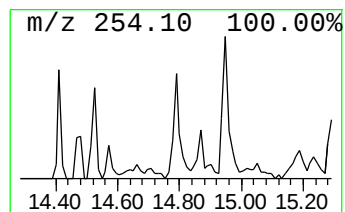
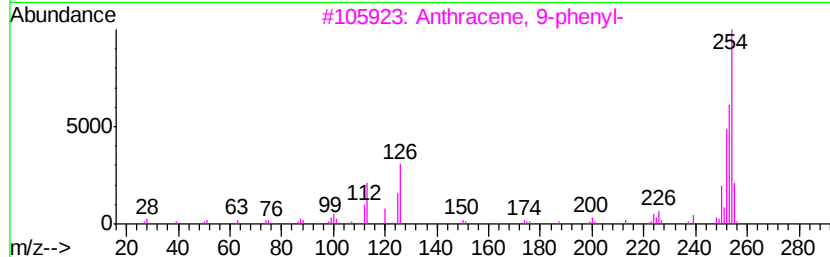
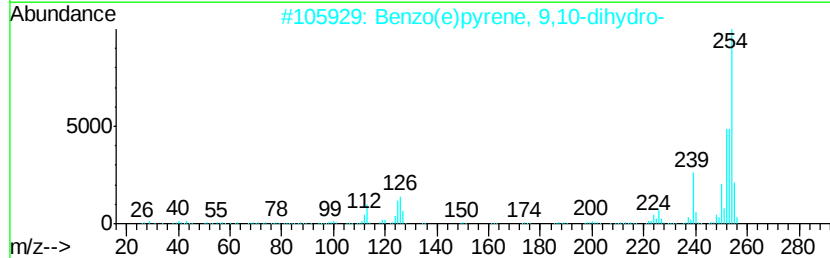
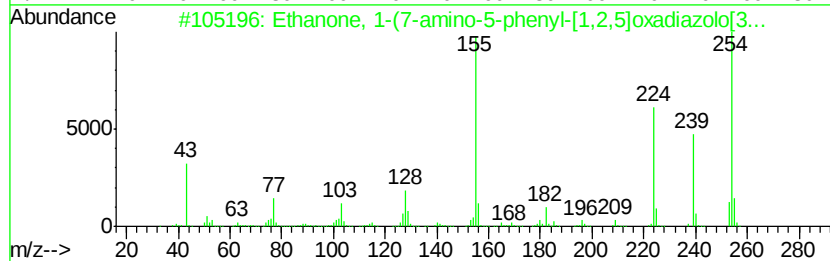
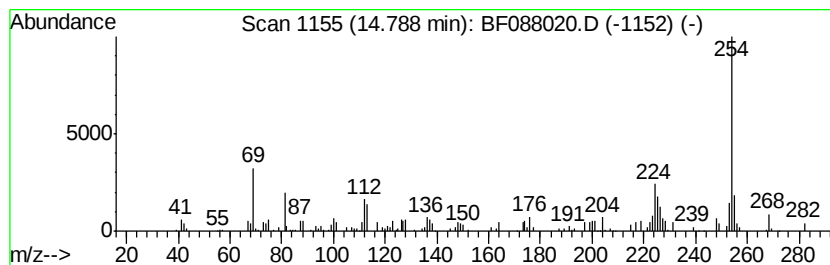
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Ethanone, 1-(7-amino-5-phen... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.79	2.94 ng	91644	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanone, 1-(7-amino-5-phenyl-[1...	254	C13H10N4O2	1000316-41-1	59
2		Benzo(e)pyrene, 9,10-dihydro-	254	C20H14	066788-01-0	58
3		Anthracene, 9-phenyl-	254	C20H14	000602-55-1	53
4		3-Hydroxy-1-methoxyvanthraquinone	254	C15H10O4	028504-24-7	53
5		3-(2-Furan-2-yl-2-oxo-ethylidene...	254	C14H10N2O3	1000297-32-1	53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF061416\
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Acq On : 15 Jun 2016 1:50
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ALS Vial : 24 Sample Multiplier: 1

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ClientSampleId :
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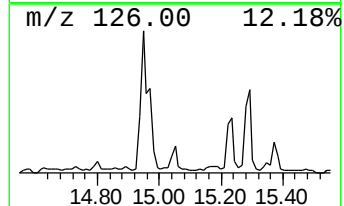
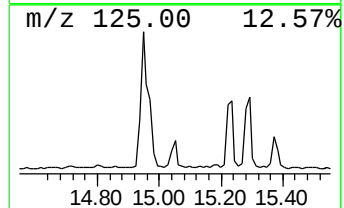
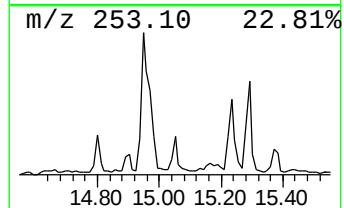
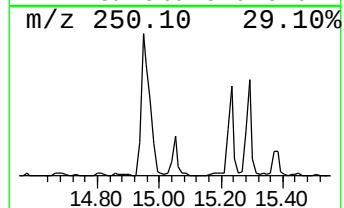
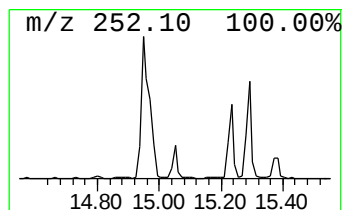
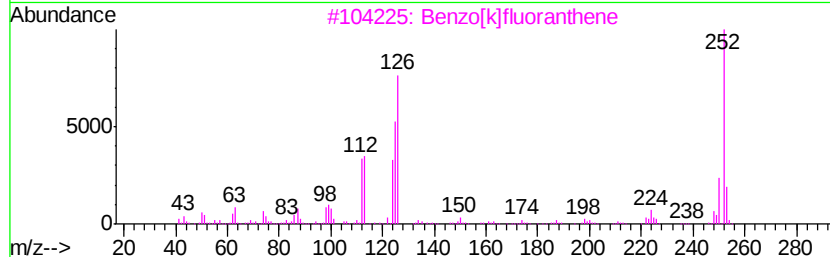
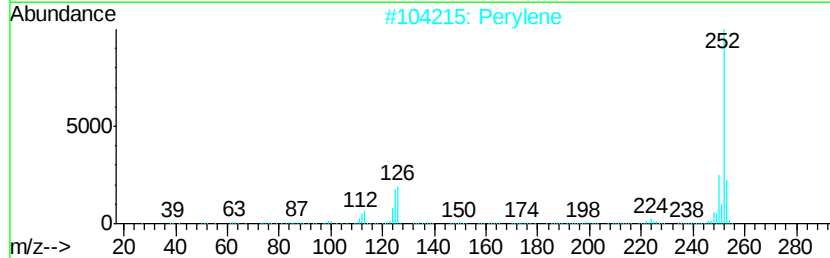
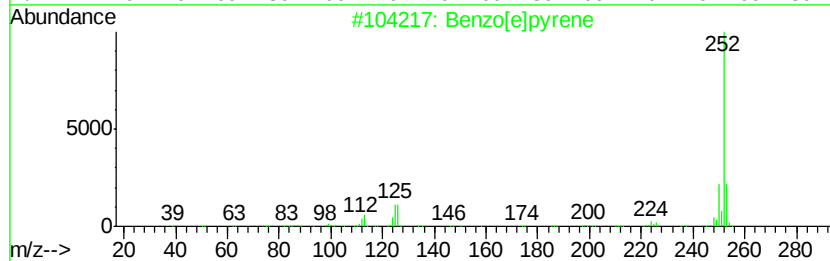
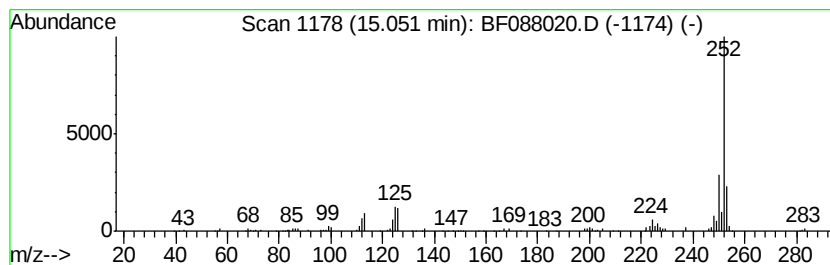
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Benzo[e]pyrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.05	3.29 ng	102825	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	96
2		Perylene	252	C20H12	000198-55-0	95
3		Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
4		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	93
5		Benzo[a]pyrene	252	C20H12	000050-32-8	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061416\
Data File : BF088020.D
Acq On : 15 Jun 2016 1:50
Operator : UM/SJ
Sample : H3471-07
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
STA-1037-(0-4)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.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
Propane, 1,1-dime...	2.00	13.9	ng	389084	1	6.79	559864	20.0
2-Pentanone, 4-hy...	4.96	6.8	ng	190579	1	6.79	559864	20.0
unknown6.55	6.55	89.4	ng	2501660	1	6.79	559864	20.0
Ethanol, 2-(2-eth...	6.62	3.6	ng	101955	1	6.79	559864	20.0
Tridecane	10.73	2.7	ng	121388	4	11.30	890771	20.0
4H-Cyclopenta[def...	11.92	2.5	ng	111536	4	11.30	890771	20.0
11H-Benzo[b]fluorene	13.07	2.1	ng	161531	5	13.94	1505880	20.0
Pentafluoropropio...	13.79	2.1	ng	156976	5	13.94	1505880	20.0
1-Trifluorosilylt...	14.72	2.3	ng	70268	6	15.35	624285	20.0
Ethanone, 1-(7-am...	14.79	2.9	ng	91644	6	15.35	624285	20.0
Benzo[e]pyrene	15.05	3.3	ng	102825	6	15.35	624285	20.0