

Data Path : Z:\HPCHEM1\BNA F\DATA\BF063016\
 Data File : BF088561.D
 Acq On : 1 Jul 2016 6:38
 Operator : UM/SJ
 Sample : H3747-08
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C8-B4(0-4.5)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-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	1.690	5	9	11	rVB	5001308	7159732	100.00%	16.433%
2	1.735	12	13	23	rVV	331182	684009	9.55%	1.570%
3	1.872	23	25	41	rVV	2092092	3586351	50.09%	8.231%
4	2.078	41	43	58	rVB	106538	280170	3.91%	0.643%
5	3.084	128	131	137	rBV	51392	110236	1.54%	0.253%
6	5.004	296	299	303	rBV	178931	246132	3.44%	0.565%
7	5.416	332	335	338	rBV	2497116	2909584	40.64%	6.678%
8	6.456	422	426	428	rBV	2770363	3567168	49.82%	8.187%
9	6.581	434	437	440	rBV	2450990	3098176	43.27%	7.111%
10	6.822	455	458	460	rBV	855363	886719	12.38%	2.035%
11	6.970	469	471	474	rVB	1900925	2499015	34.90%	5.736%
12	7.382	504	507	509	rBV	2072958	2050729	28.64%	4.707%
13	8.102	567	570	572	rBV	1417104	1189751	16.62%	2.731%
14	9.176	661	664	667	rBV	2727482	3331363	46.53%	7.646%
15	9.565	697	698	701	rBV	126517	165023	2.30%	0.379%
16	9.850	721	723	725	rBV	1285080	1179887	16.48%	2.708%
17	10.639	789	792	794	rBV	1688942	1712094	23.91%	3.929%
18	11.211	840	842	848	rVB	73375	84590	1.18%	0.194%
19	11.325	850	852	856	rVV	780882	1101097	15.38%	2.527%
20	11.862	897	899	901	rBV	93057	126810	1.77%	0.291%
21	12.022	911	913	920	rVB2	38033	71609	1.00%	0.164%
22	12.536	956	958	960	rBV	285971	243636	3.40%	0.559%
23	12.754	974	977	979	rBV2	320470	543941	7.60%	1.248%
24	12.822	979	983	985	rVB3	95011	125714	1.76%	0.289%
25	12.914	989	991	994	rVB	2309917	2214838	30.93%	5.083%
26	13.154	1009	1012	1014	rBV	63007	86268	1.20%	0.198%
27	13.451	1036	1038	1040	rVV	191724	174405	2.44%	0.400%
28	13.496	1040	1042	1047	rVB2	73926	121971	1.70%	0.280%
29	13.702	1056	1060	1062	rBV3	35246	74250	1.04%	0.170%
30	13.828	1069	1071	1074	rBV	216533	223443	3.12%	0.513%
31	13.954	1079	1082	1087	rBV2	923768	1129934	15.78%	2.593%
32	14.148	1096	1099	1100	rBV	70590	103473	1.45%	0.237%
33	14.445	1123	1125	1131	rBV3	109438	258580	3.61%	0.593%
34	14.742	1149	1151	1153	rBV	88380	76191	1.06%	0.175%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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35	14.879	1161	1163	1165	rBV	141598	158324	2.21%	0.363%
36	14.959	1168	1170	1175	rVB	130061	263829	3.68%	0.606%
37	15.062	1177	1179	1180	rVV	132304	135507	1.89%	0.311%
38	15.097	1180	1182	1187	rVB2	115724	183197	2.56%	0.420%
39	15.245	1193	1195	1197	rVV	67685	75908	1.06%	0.174%
40	15.302	1197	1200	1202	rVV	94449	137548	1.92%	0.316%
41	15.359	1202	1205	1208	rVV	614340	837974	11.70%	1.923%
42	15.417	1209	1210	1216	rVB	77887	105174	1.47%	0.241%
43	15.600	1224	1226	1230	rBV2	53253	102114	1.43%	0.234%
44	15.828	1243	1246	1248	rBV	91594	153965	2.15%	0.353%

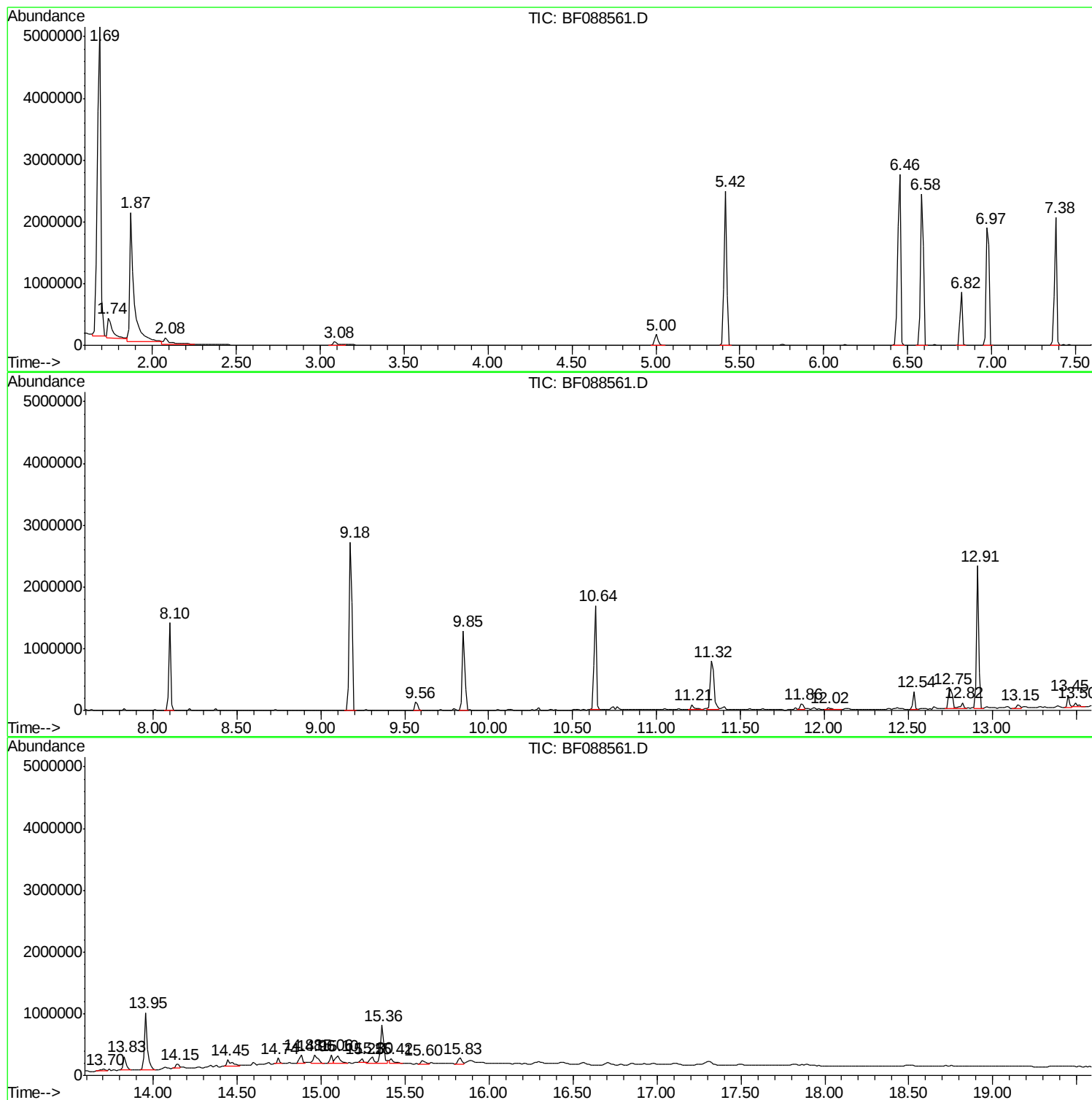
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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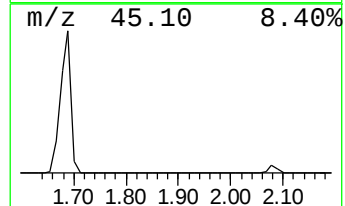
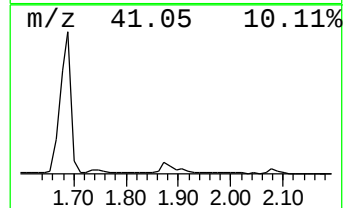
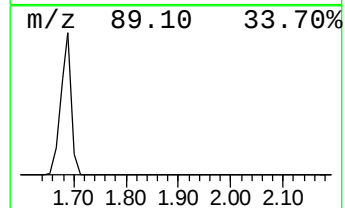
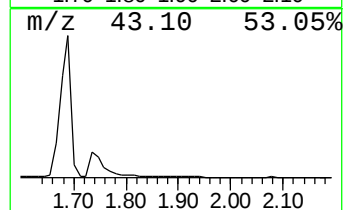
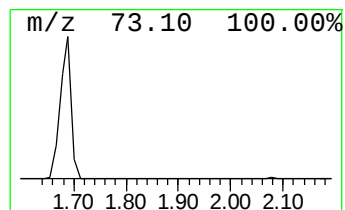
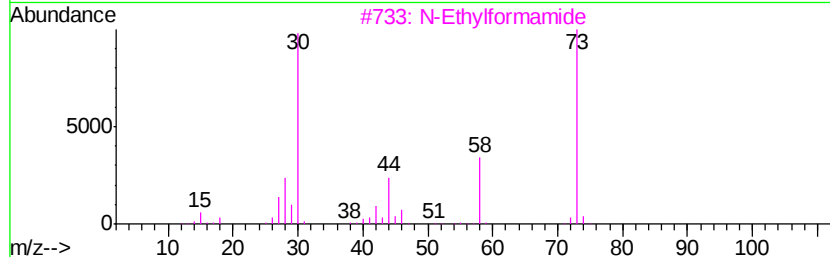
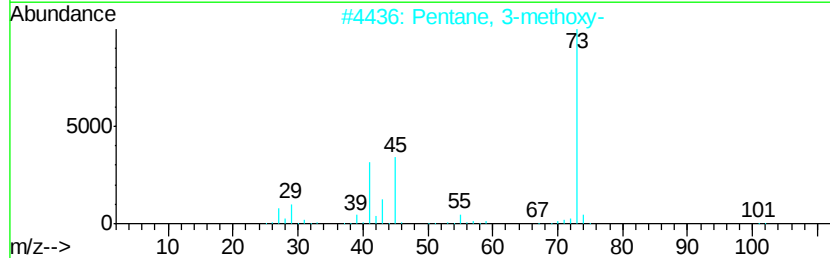
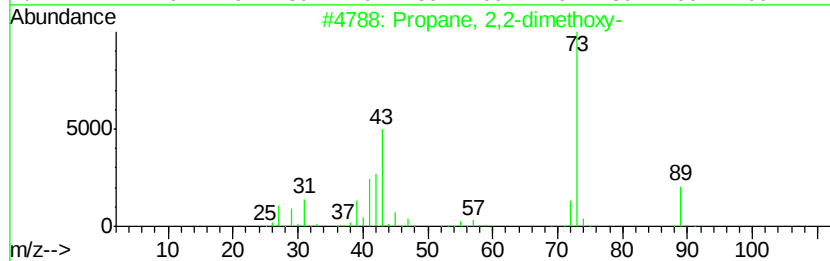
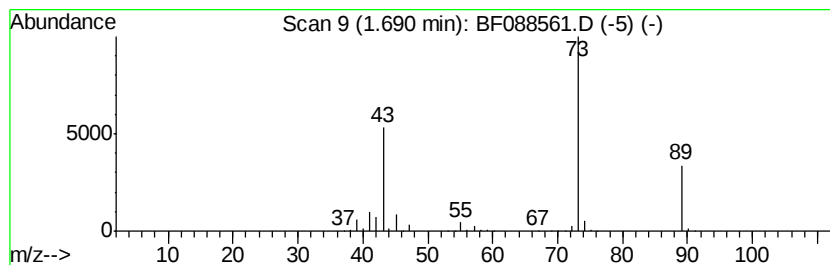
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.69	161.49 ng	7159730	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	56
2		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
3		N-Ethylformamide	73	C3H7NO	000627-45-2	9
4		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9
5		1,3-Diaminoguanidine	89	CH7N5	004364-78-7	9



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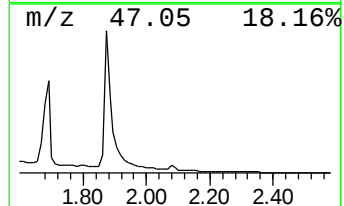
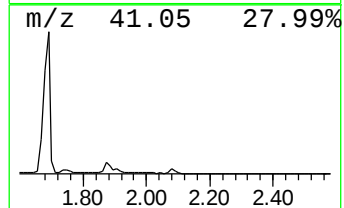
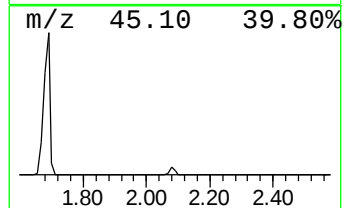
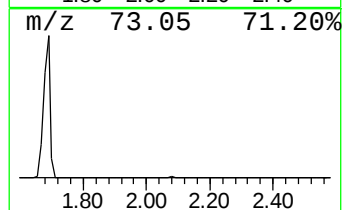
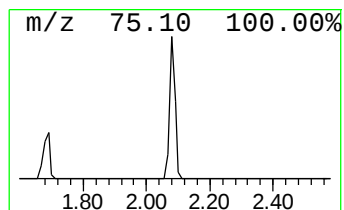
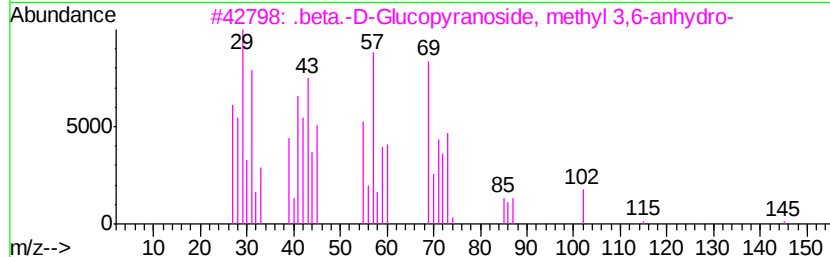
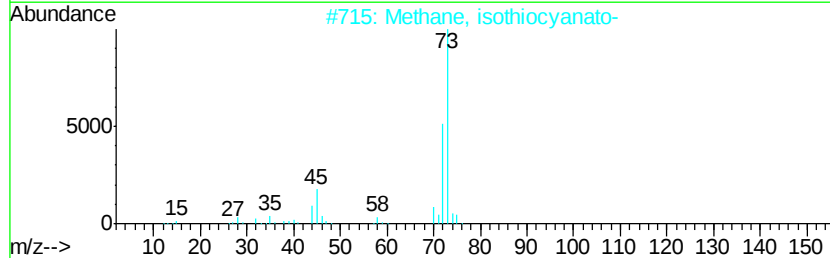
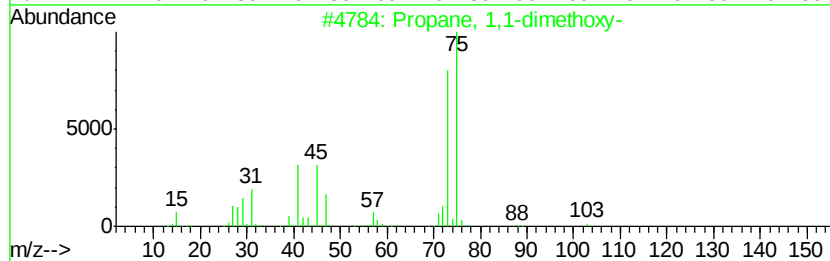
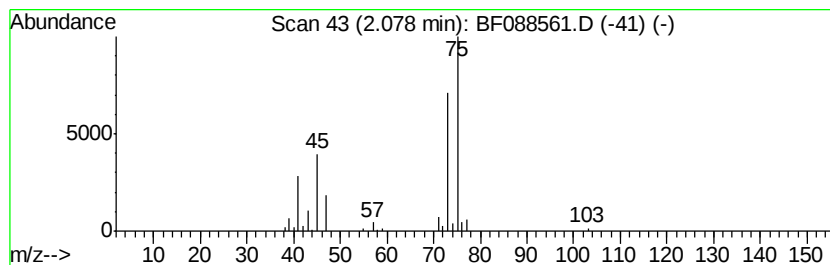
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Peak Number 4 Propane, 1,1-dimethoxy- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.08	6.32 ng	280170	1,4-Dichlorobenzene-d4	6.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	64
2			Methane, isothiocyanato-	73	C2H3NS	000556-61-6	9
3			.beta.-D-Glucopyranoside, methyl...	176	C7H12O5	003056-46-0	9
4			Ethane, 2-chloro-1,1-dimethoxy-	124	C4H9ClO2	000097-97-2	9
5			Propane, 1,1-dimethoxy-2-methyl-	118	C6H14O2	041632-89-7	9



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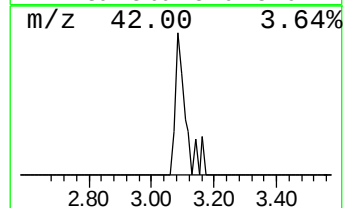
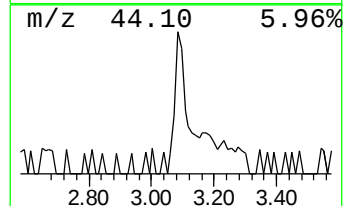
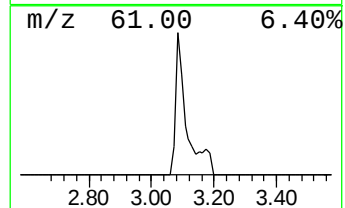
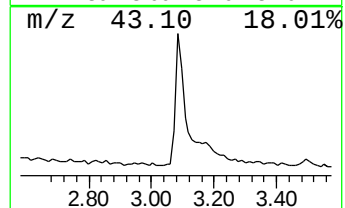
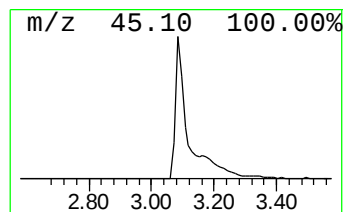
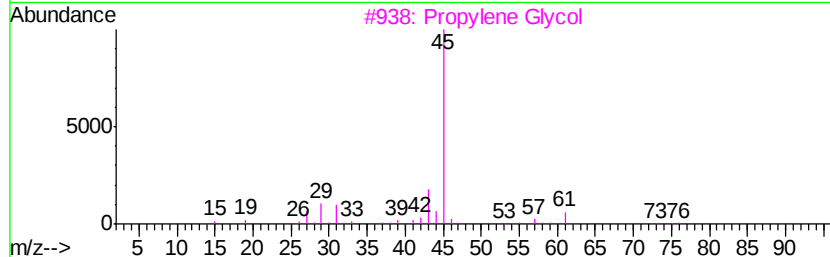
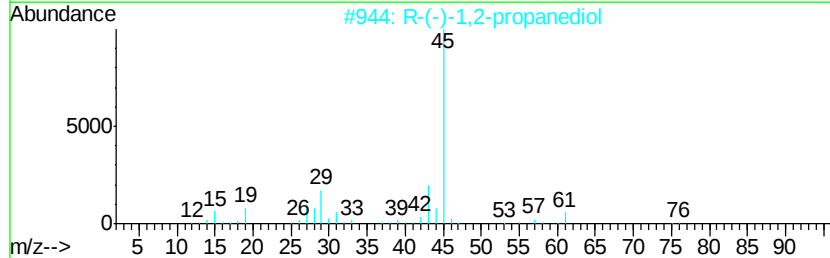
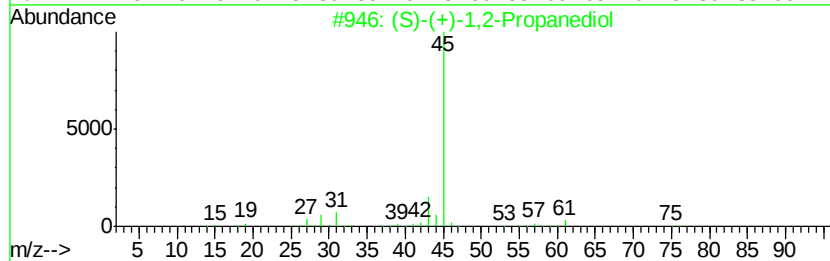
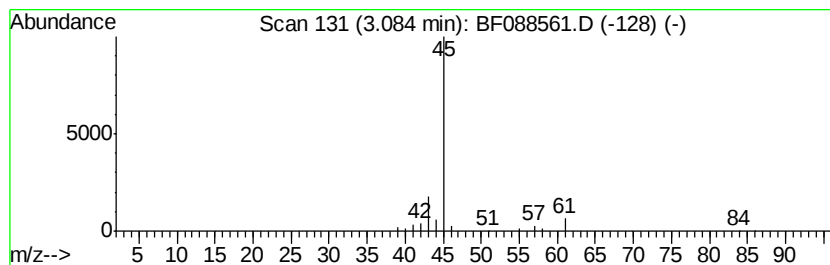
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (S)-(+)-1,2-Propanediol Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.08	2.49 ng	110236	1,4-Dichlorobenzene-d4	6.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	64
2	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	43
3	Propylene Glycol	76	C3H8O2	000057-55-6	9
4	Isopropyl Alcohol	60	C3H8O	000067-63-0	9
5	2-Heptanol	116	C7H16O	000543-49-7	9



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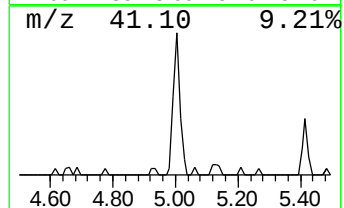
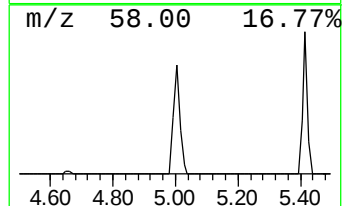
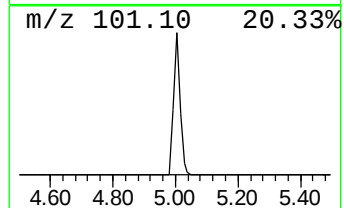
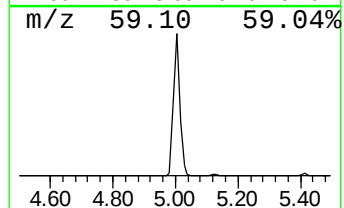
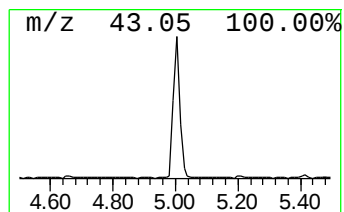
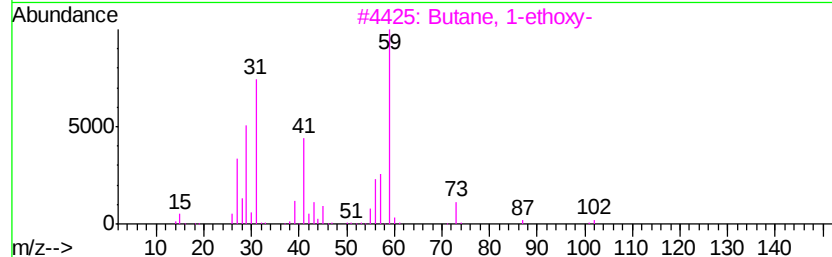
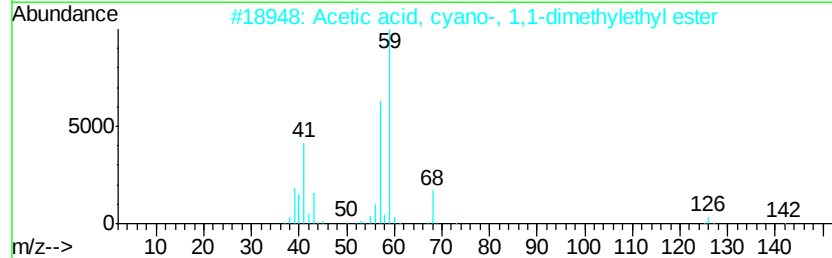
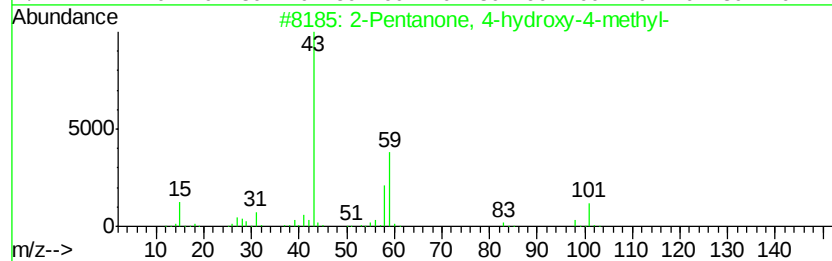
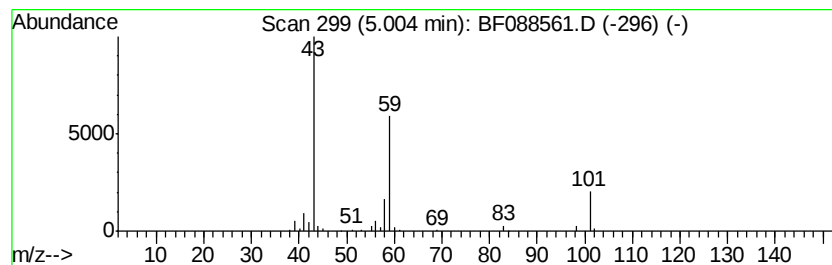
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Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.00	5.55 ng	246132	1,4-Dichlorobenzene-d4	6.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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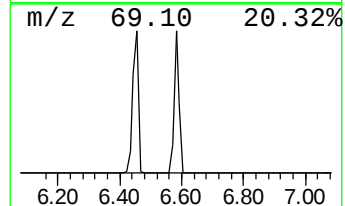
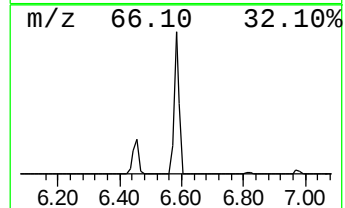
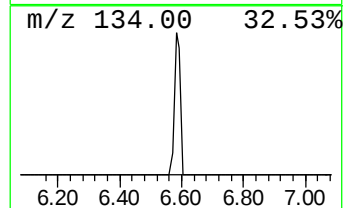
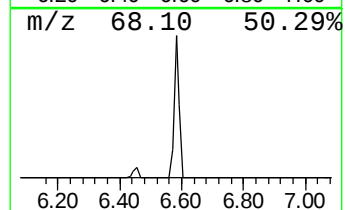
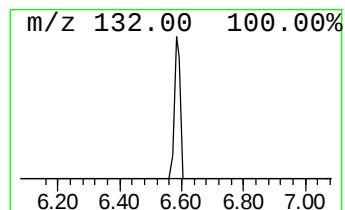
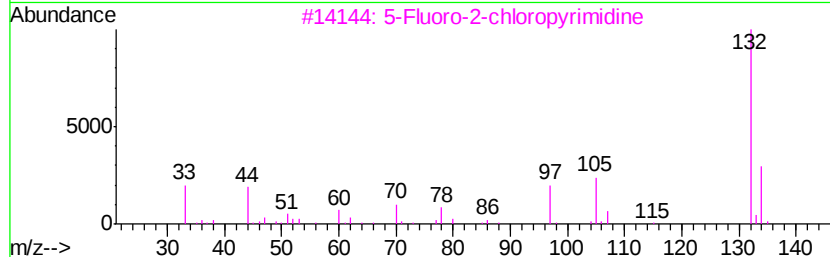
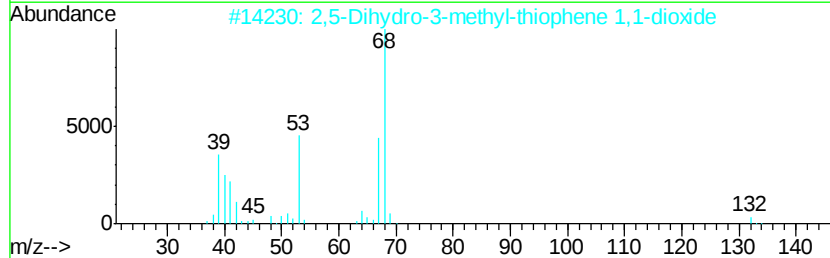
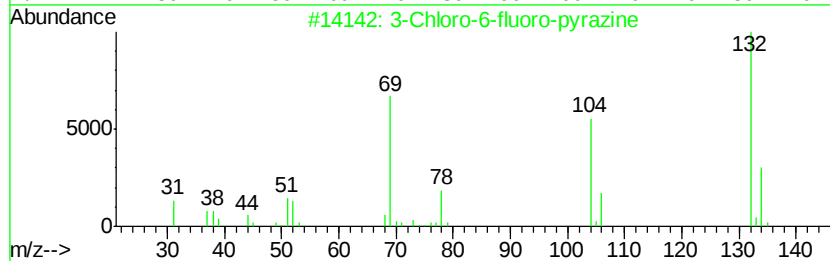
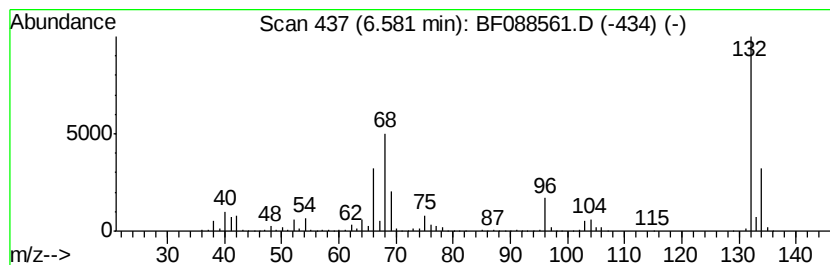
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown6.58 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	69.88 ng	3098180	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		2,5-Dihydro-3-methyl-thiophene 1...	132	C5H8O2S	001193-10-8	17
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
5		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
Data File : BF088561.D
Acq On : 1 Jul 2016 6:38
Operator : UM/SJ
Sample : H3747-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

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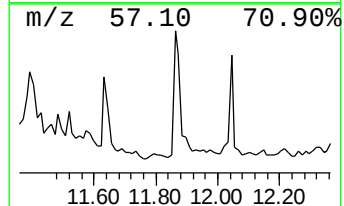
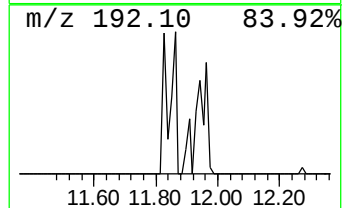
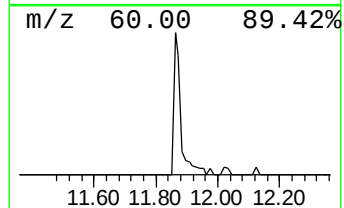
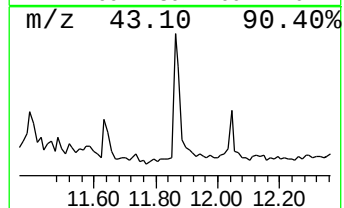
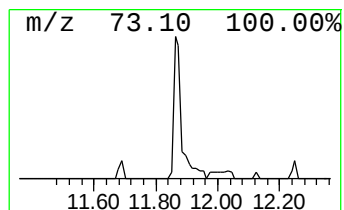
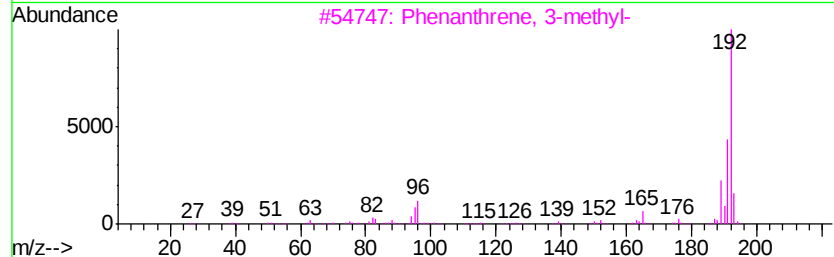
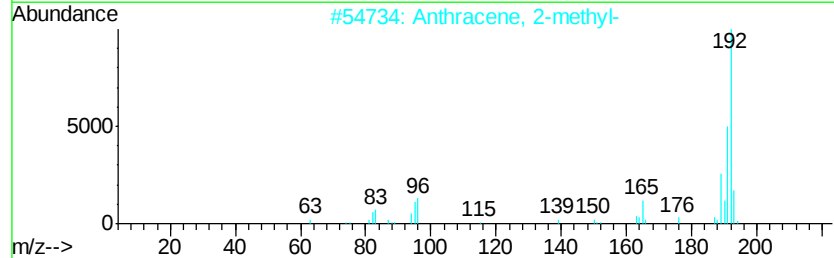
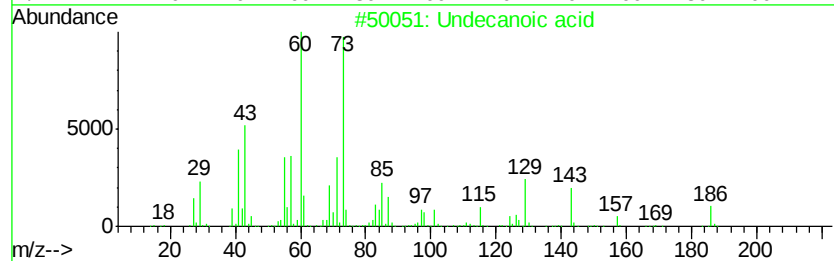
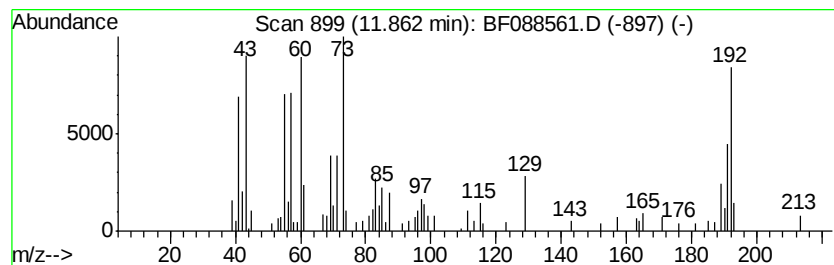
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown11.86 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	2.30 ng	126810	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Undecanoic acid	186	C11H22O2	000112-37-8	46
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	41
3		Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	35
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	35
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	35



Data Path : Z:\HPCHEM1\BNA F\DATA\BF063016\
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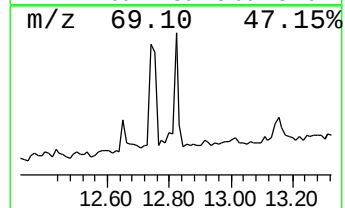
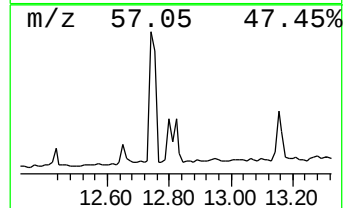
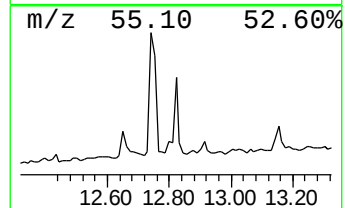
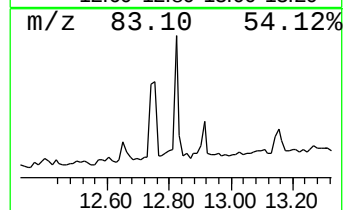
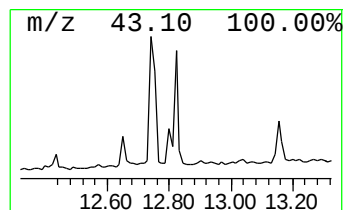
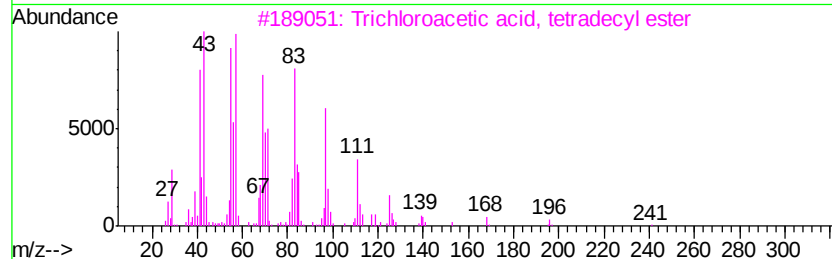
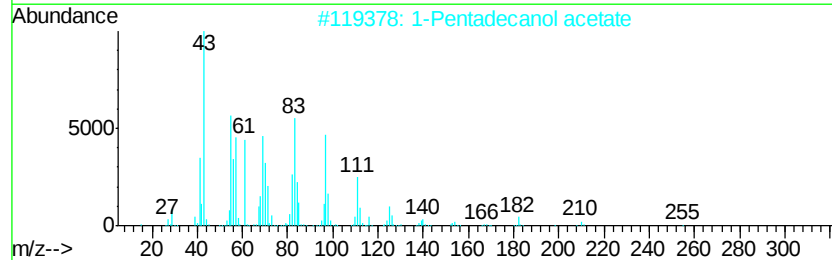
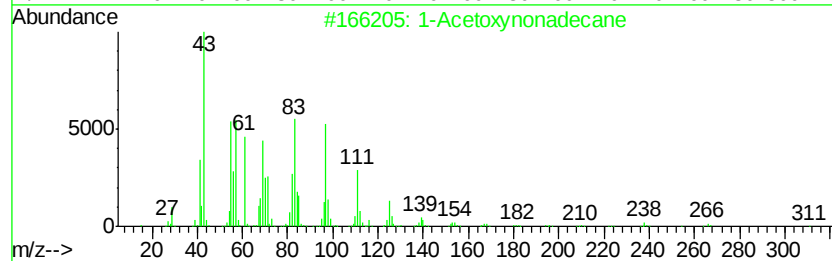
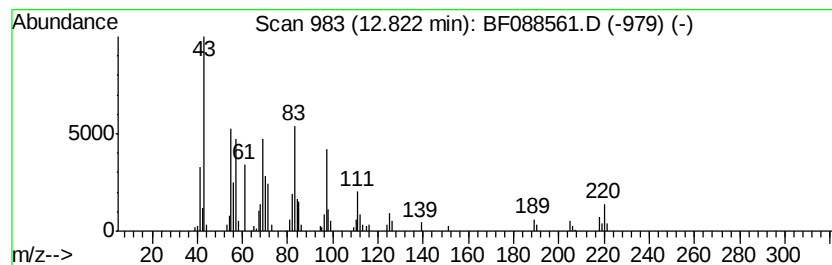
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1-Acetoxynonadecane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.82	2.23 ng	125714	Chrysene-d12	13.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Acetoxynonadecane	326	C21H42O2	001577-43-1	91
2			1-Pentadecanol acetate	270	C17H34O2	000629-58-3	90
3			Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	87
4			Eicosyl acetate	340	C22H44O2	1000351-77-9	87
5			Behenic alcohol	326	C22H46O	000661-19-8	87



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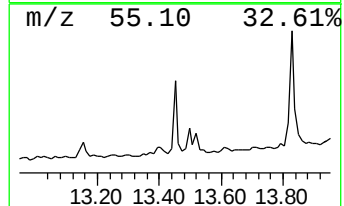
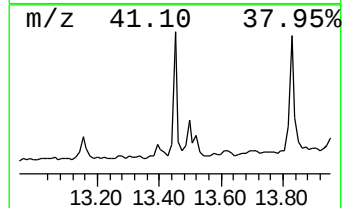
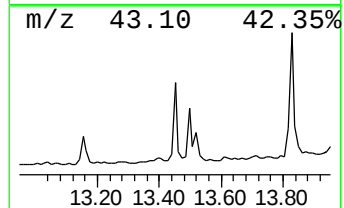
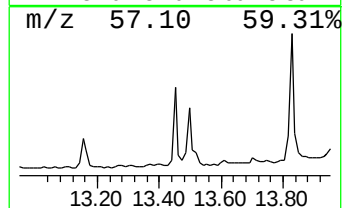
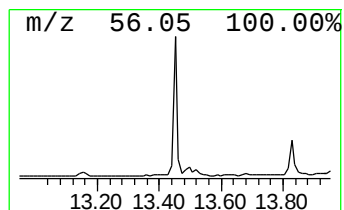
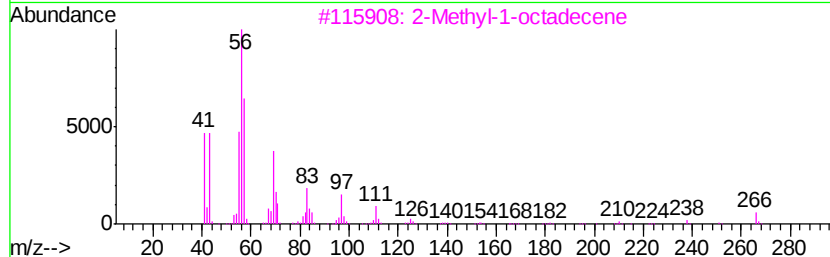
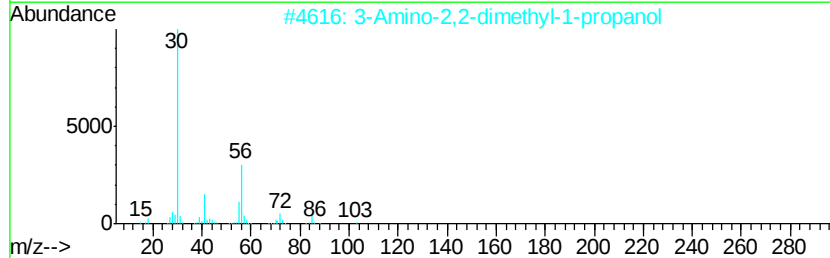
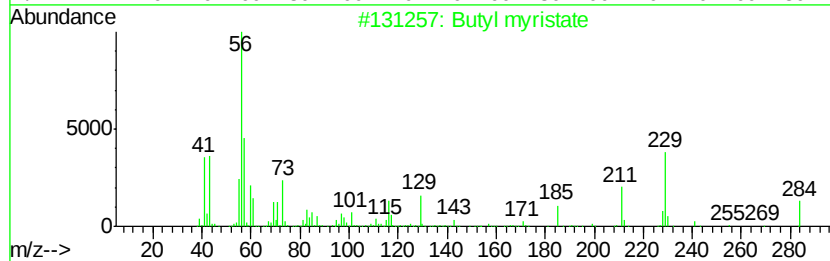
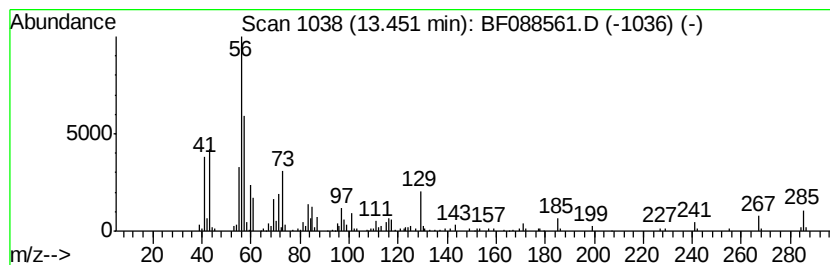
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Butyl myristate Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.45	3.09 ng	174405	Chrysene-d12	13.95	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl myristate	284	C18H36O2	000110-36-1	86
2	3-Amino-2,2-dimethyl-1-propanol	103	C5H13NO	026734-09-8	38
3	2-Methyl-1-octadecene	266	C19H38	061868-20-0	38
4	Urea, N,N'-di-2-propenyl-	140	C7H12N2O	001801-72-5	35
5	Cyclopropane, 1-(1-methylethyl)-...	210	C15H30	041977-39-3	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
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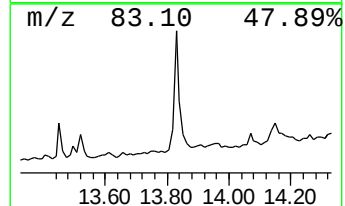
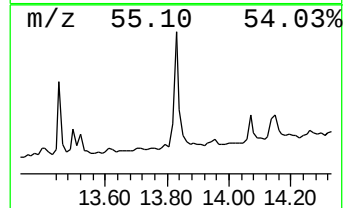
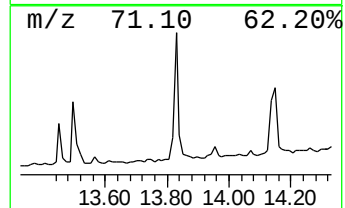
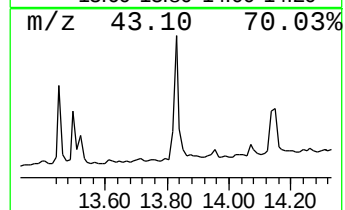
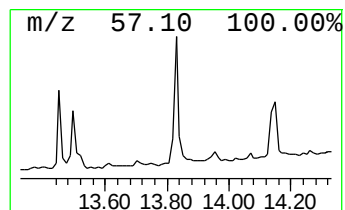
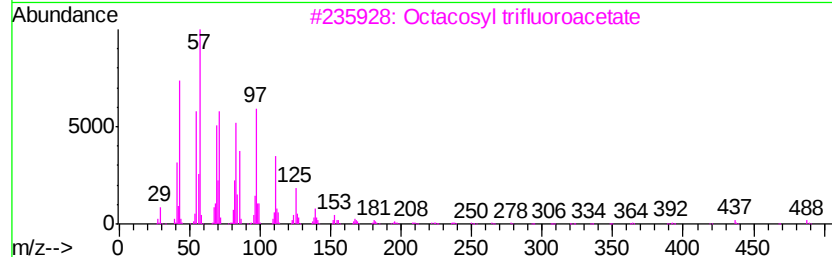
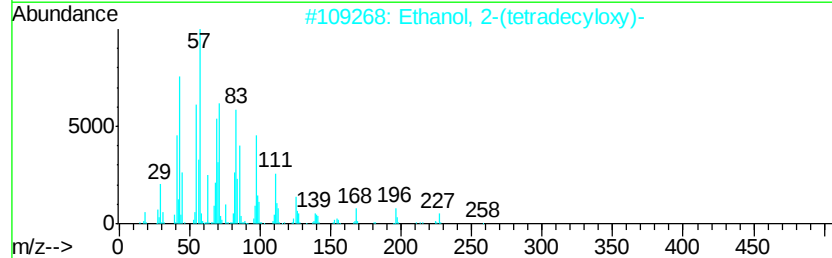
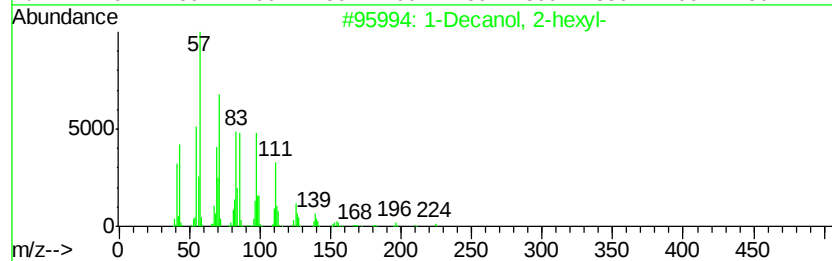
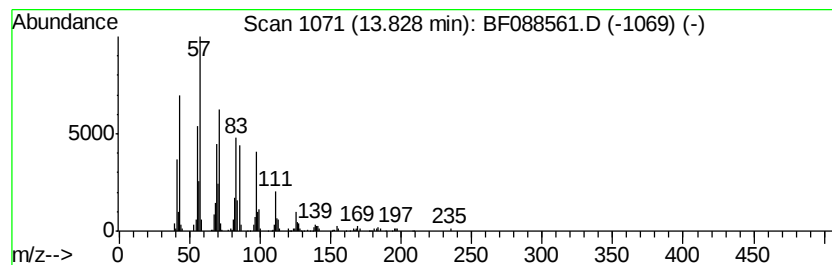
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1-Decanol, 2-hexyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	3.95 ng	223443	Chrysene-d12	13.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	91
2			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	87
3			Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	87
4			Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	76
5			Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	76



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
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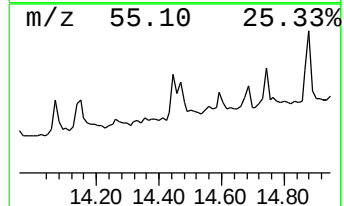
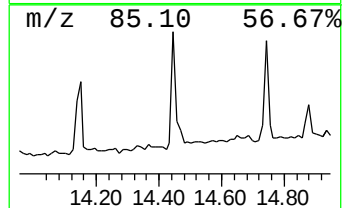
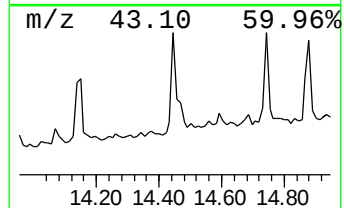
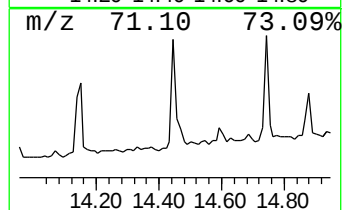
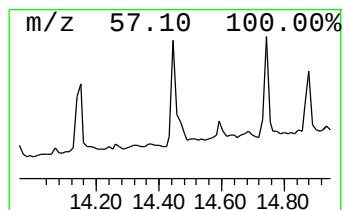
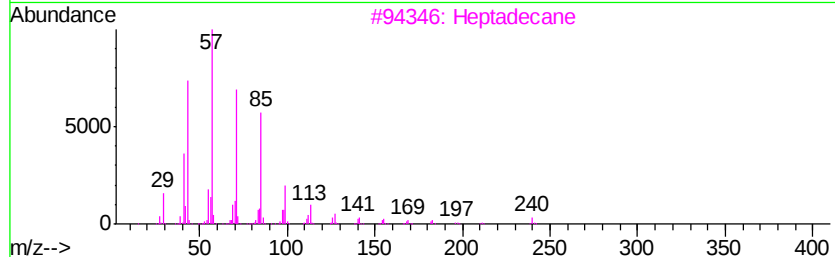
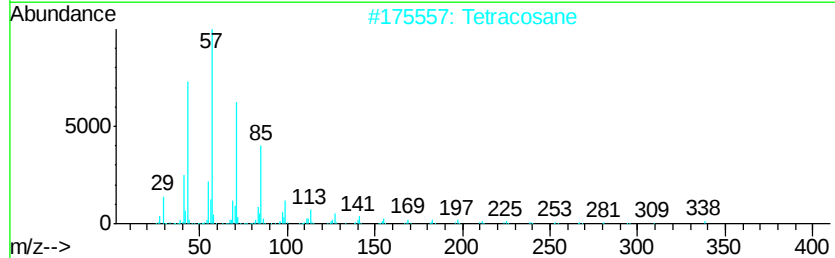
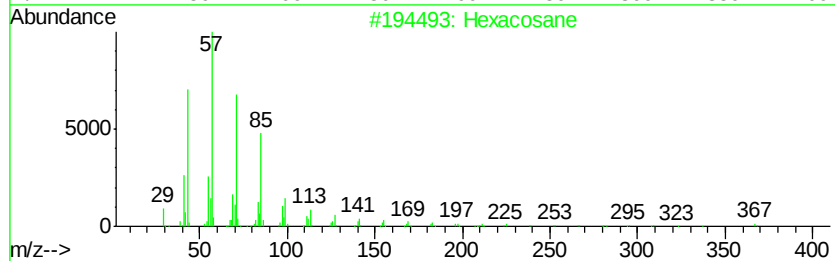
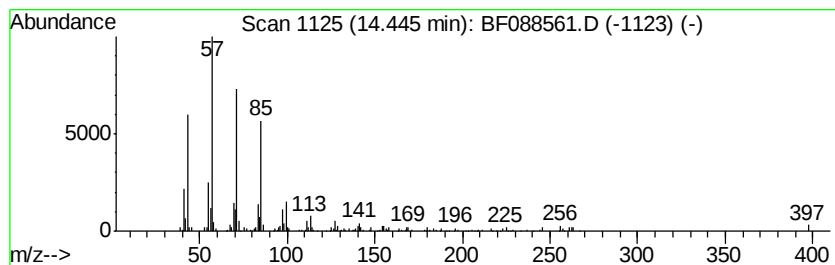
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Hexacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.45	4.58 ng	258580	Chrysene-d12	13.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	87
2			Tetracosane	338	C24H50	000646-31-1	87
3			Heptadecane	240	C17H36	000629-78-7	86
4			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	83
5			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
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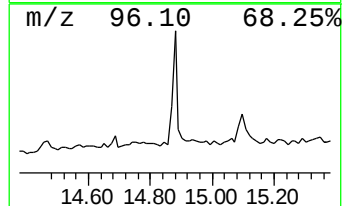
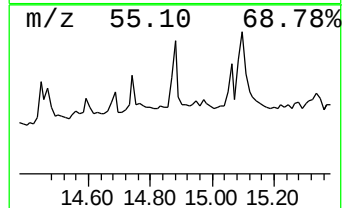
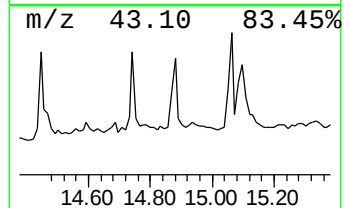
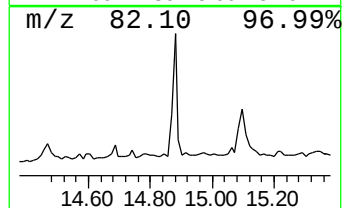
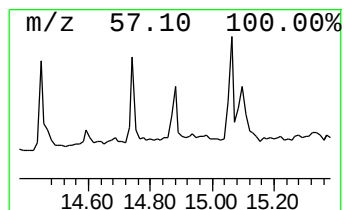
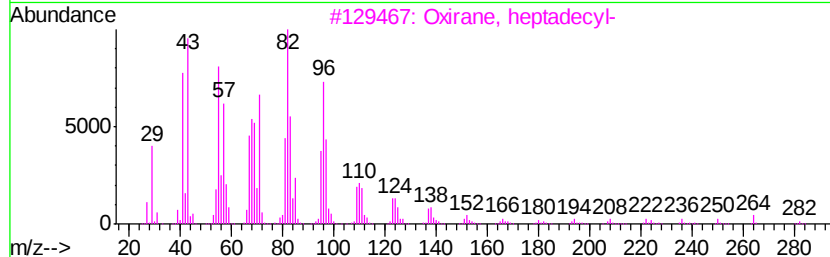
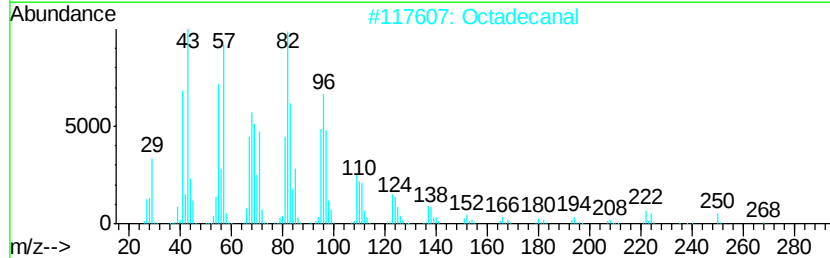
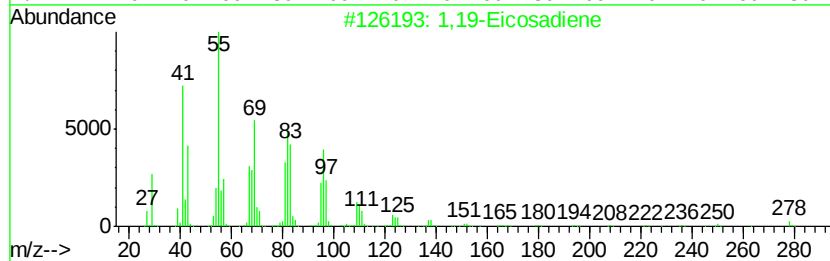
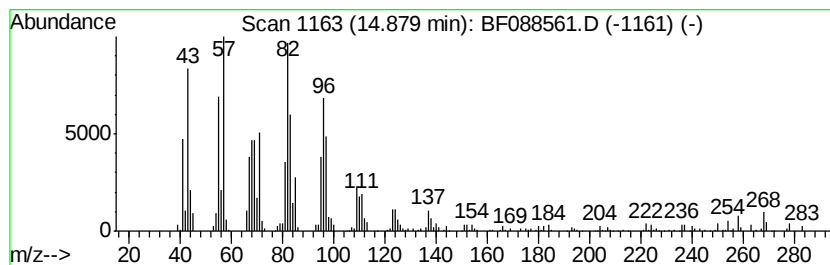
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 1,19-Eicosadiene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.88	3.78 ng	158324	Perylene-d12	15.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,19-Eicosadiene	278	C20H38	014811-95-1	92
2	Octadecanal	268	C18H36O	000638-66-4	91
3	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	90
4	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	87
5	Z-2-Dodecenol	184	C12H24O	069064-36-4	76



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
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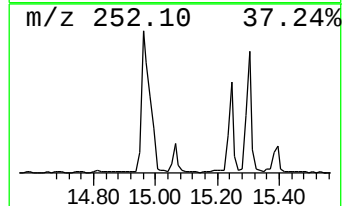
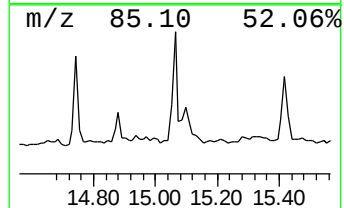
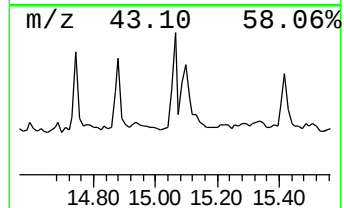
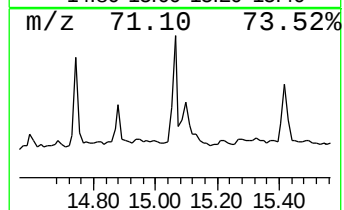
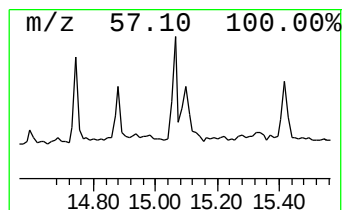
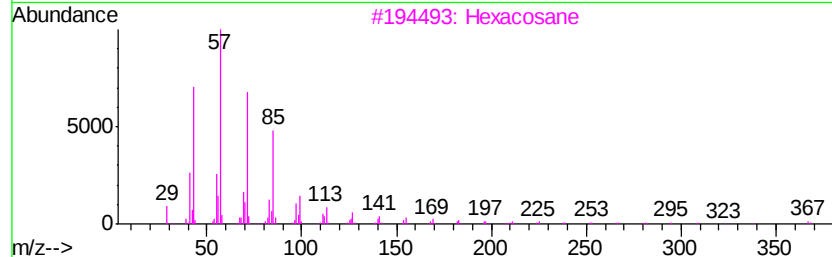
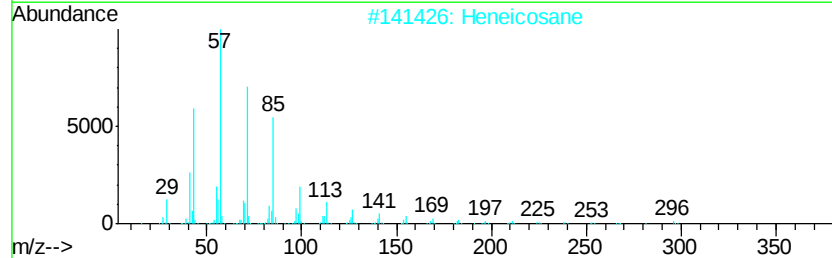
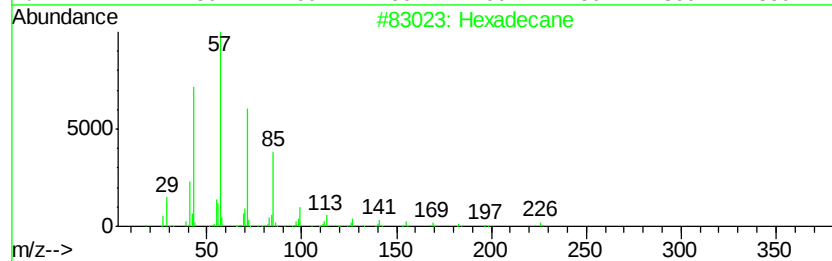
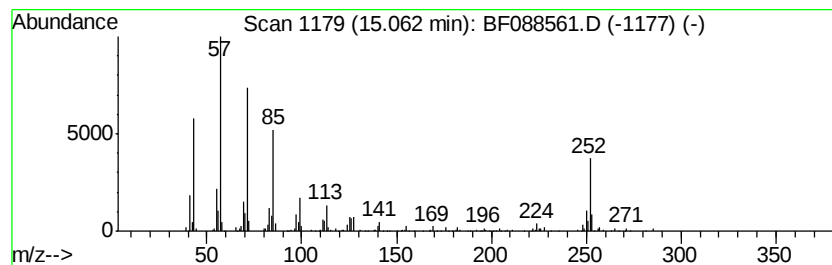
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BNA_F
ClientSampleId :
C8-B4(0-4.5)C

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 16 Hexadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.06	3.23 ng	135507	Perylene-d12		15.36
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	76
2	Heneicosane	296	C21H44	000629-94-7	76
3	Hexacosane	366	C26H54	000630-01-3	76
4	Tetracosane	338	C24H50	000646-31-1	64
5	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
Data File : BF088561.D
Acq On : 1 Jul 2016 6:38
Operator : UM/SJ
Sample : H3747-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
C8-B4(0-4.5)C

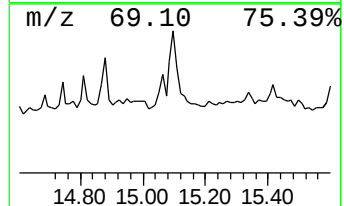
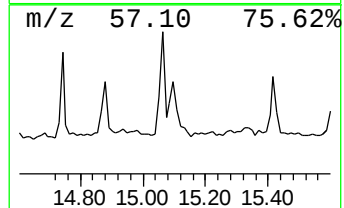
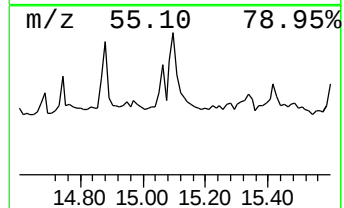
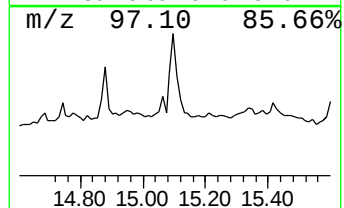
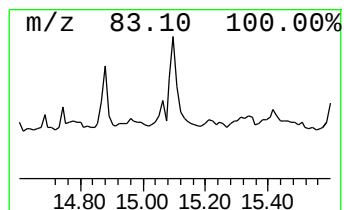
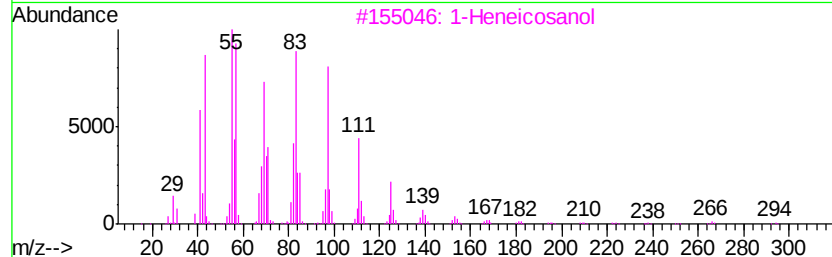
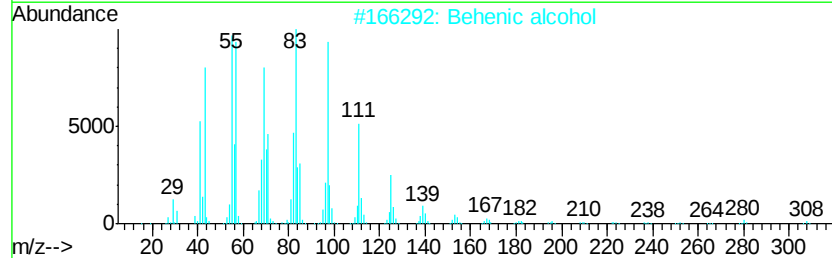
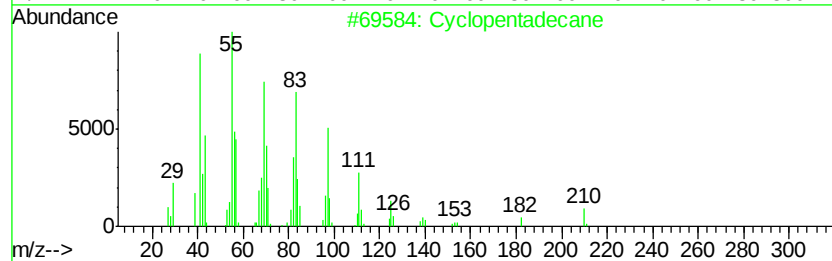
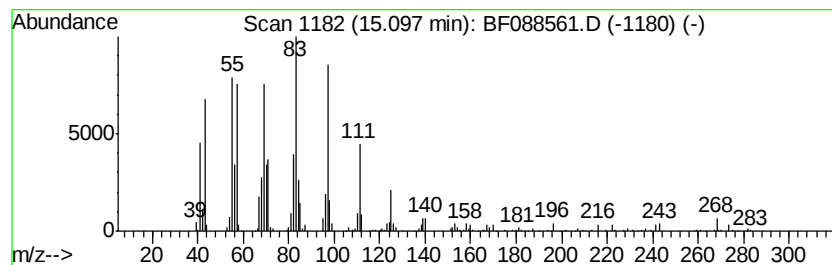
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Cyclopentadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.10	4.37 ng	183197	Perylene-d12	15.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentadecane	210	C15H30	000295-48-7	93
2		Behenic alcohol	326	C22H46O	000661-19-8	91
3		1-Heneicosanol	312	C21H44O	015594-90-8	91
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	87
5		Cyclotetradecane	196	C14H28	000295-17-0	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
Data File : BF088561.D
Acq On : 1 Jul 2016 6:38
Operator : UM/SJ
Sample : H3747-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

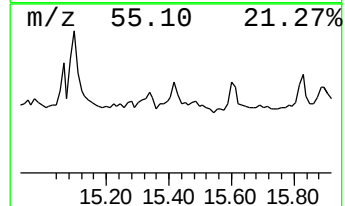
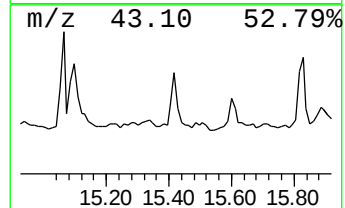
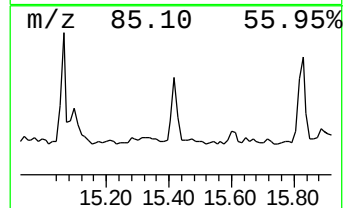
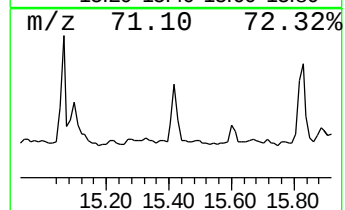
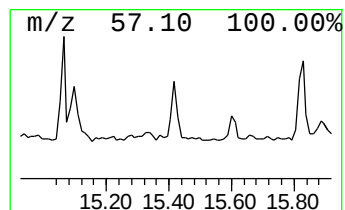
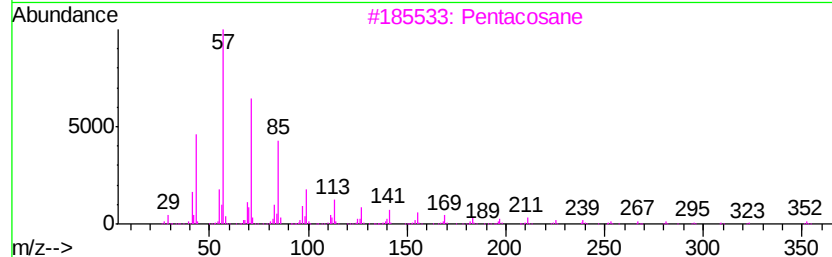
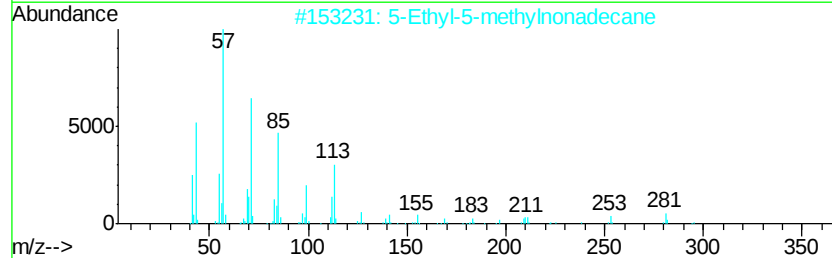
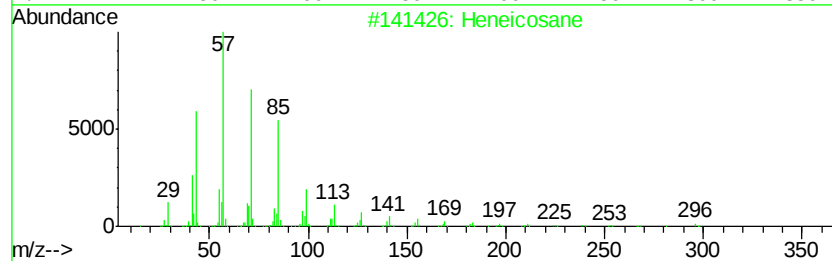
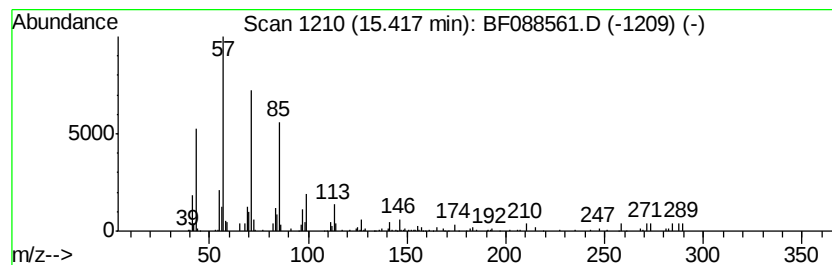
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Heneicosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.42	2.51 ng	105174	Perylene-d12	15.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	76
2	5-Ethyl-5-methylnonadecane	310	C22H46	1000360-42-7	74
3	Pentacosane	352	C25H52	000629-99-2	68
4	1-Iodoundecane	282	C11H23I	004282-44-4	68
5	Nonadecane	268	C19H40	000629-92-5	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF063016\
Data File : BF088561.D
Acq On : 1 Jul 2016 6:38
Operator : UM/SJ
Sample : H3747-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

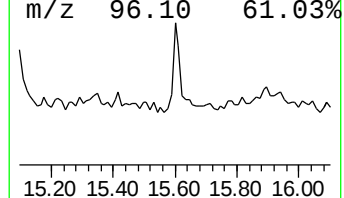
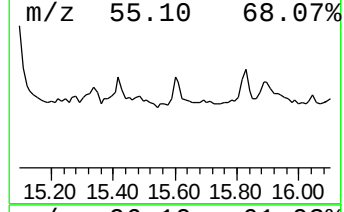
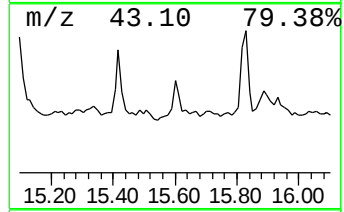
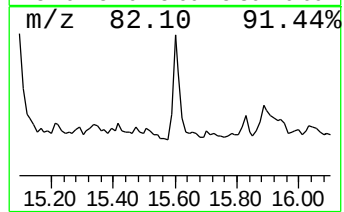
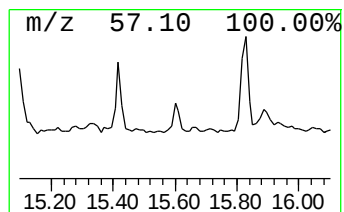
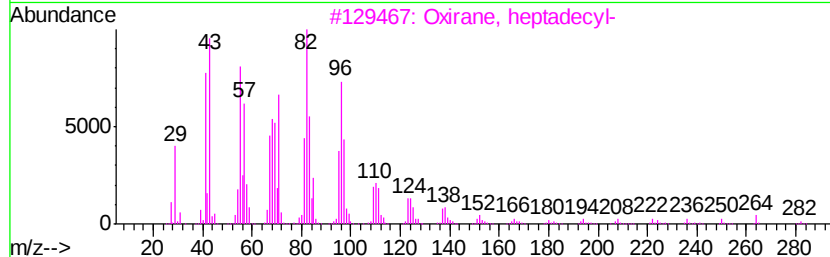
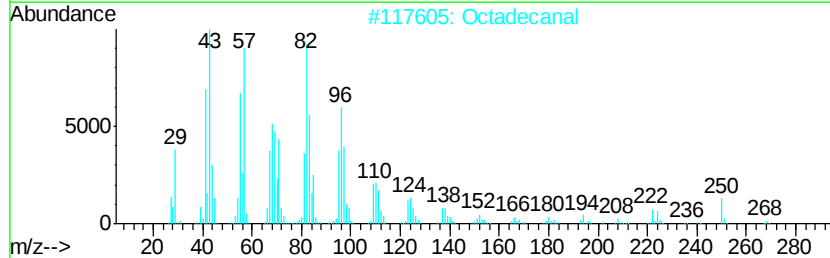
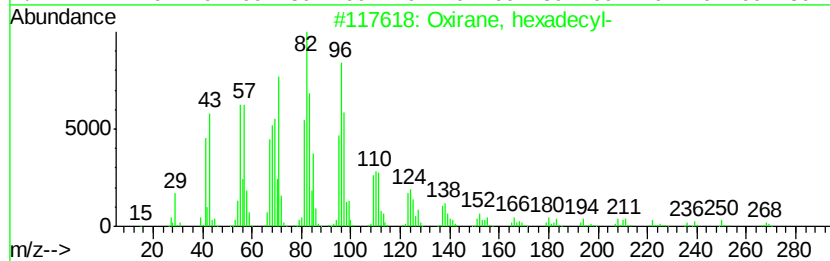
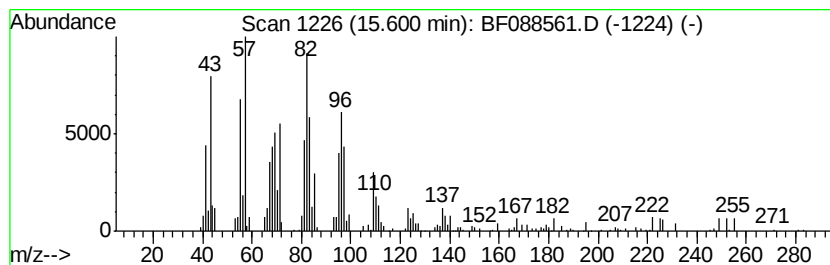
Instrument :
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ClientSampleId :
C8-B4(0-4.5)C

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 19 Oxirane, hexadecyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.60	2.44 ng	102114	Perylene-d12	15.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	90
2	Octadecanal	268	C18H36O	000638-66-4	86
3	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	78
4	Pentadecanal-	226	C15H30O	002765-11-9	72
5	1,19-Eicosadiene	278	C20H38	014811-95-1	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
 Data File : BF088561.D
 Acq On : 1 Jul 2016 6:38
 Operator : UM/SJ
 Sample : H3747-08
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C8-B4(0-4.5)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
Propane, 2,2-dime...	1.69	161.5	ng	7159730	1	6.82	886719	20.0
Propane, 1,1-dime...	2.08	6.3	ng	280170	1	6.82	886719	20.0
(S)-(+)-1,2-Propa...	3.08	2.5	ng	110236	1	6.82	886719	20.0
2-Pentanone, 4-hy...	5.00	5.5	ng	246132	1	6.82	886719	20.0
unknown6.58	6.58	69.9	ng	3098180	1	6.82	886719	20.0
unknown11.86	11.86	2.3	ng	126810	4	11.32	1101100	20.0
1-Acetoxynonadecane	12.82	2.2	ng	125714	5	13.95	1129930	20.0
Butyl myristate	13.45	3.1	ng	174405	5	13.95	1129930	20.0
1-Decanol, 2-hexyl-	13.83	4.0	ng	223443	5	13.95	1129930	20.0
Hexacosane	14.45	4.6	ng	258580	5	13.95	1129930	20.0
1,19-Eicosadiene	14.88	3.8	ng	158324	6	15.36	837974	20.0
Hexadecane	15.06	3.2	ng	135507	6	15.36	837974	20.0
Cyclopentadecane	15.10	4.4	ng	183197	6	15.36	837974	20.0
Heneicosane	15.42	2.5	ng	105174	6	15.36	837974	20.0
Oxirane, hexadecyl-	15.60	2.4	ng	102114	6	15.36	837974	20.0