

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
 Data File : BF088551.D
 Acq On : 1 Jul 2016 1:48
 Operator : UM/SJ
 Sample : H3630-05
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B1409-TP02W-1224

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	12	rVB	4653359	8699461	100.00%	16.631%
2	1.735	12	13	23	rVV	343348	749822	8.62%	1.433%
3	1.873	23	25	41	rVV	1883020	3219880	37.01%	6.155%
4	2.090	41	44	50	rVB	96922	173151	1.99%	0.331%
5	5.004	296	299	305	rBV	341696	542572	6.24%	1.037%
6	5.416	332	335	338	rBV	3002800	3572203	41.06%	6.829%
7	6.262	405	409	413	rBV3	31537	90864	1.04%	0.174%
8	6.456	422	426	428	rVV	3024020	4514180	51.89%	8.630%
9	6.582	434	437	440	rBV	3207131	3931254	45.19%	7.515%
10	6.810	456	457	459	rVV	752702	859798	9.88%	1.644%
11	6.970	468	471	473	rBV	2812470	2865896	32.94%	5.479%
12	7.382	501	507	509	rBV	2525212	2925841	33.63%	5.593%
13	7.462	512	514	517	rVB3	71672	94826	1.09%	0.181%
14	7.633	523	529	530	rVB3	42111	136226	1.57%	0.260%
15	7.747	536	539	541	rVB3	132286	170771	1.96%	0.326%
16	7.850	547	548	552	rVB	426987	358331	4.12%	0.685%
17	8.102	567	570	572	rVV	1232939	1295966	14.90%	2.477%
18	8.170	574	576	578	rVV2	81088	129246	1.49%	0.247%
19	8.399	592	596	597	rBV2	94382	134106	1.54%	0.256%
20	8.536	606	608	610	rBV2	152145	189360	2.18%	0.362%
21	8.765	624	628	629	rVV4	38990	113750	1.31%	0.217%
22	8.799	629	631	633	rVB2	62140	88208	1.01%	0.169%
23	9.130	658	660	662	rBV	129943	109935	1.26%	0.210%
24	9.176	662	664	667	rVB	3455810	3602454	41.41%	6.887%
25	9.256	669	671	674	rBV2	95634	181264	2.08%	0.347%
26	9.565	696	698	701	rVB2	681193	823592	9.47%	1.574%
27	9.771	713	716	717	rVB2	98880	136235	1.57%	0.260%
28	9.793	717	718	720	rBV2	98042	96078	1.10%	0.184%
29	9.851	720	723	725	rBV	1388496	1307234	15.03%	2.499%
30	10.113	745	746	750	rBV3	47311	96242	1.11%	0.184%
31	10.296	760	762	763	rVB	91383	89288	1.03%	0.171%
32	10.513	778	781	784	rBV	90243	116317	1.34%	0.222%
33	10.639	789	792	794	rBV	1913439	2411584	27.72%	4.610%
34	10.765	801	803	808	rVB	166806	241130	2.77%	0.461%

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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

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

35	11.325	850	852	855	rVB	1234825	1054404	12.12%	2.016%
36	11.576	872	874	878	rBV	112531	177187	2.04%	0.339%
37	11.862	897	899	904	rVV3	257059	371475	4.27%	0.710%
38	12.651	966	968	970	rBV	102337	129882	1.49%	0.248%
39	12.914	988	991	993	rBV	2622789	2558085	29.41%	4.890%
40	13.828	1070	1071	1074	rBV	185949	238642	2.74%	0.456%
41	13.954	1080	1082	1084	rBV	761582	694241	7.98%	1.327%
42	15.360	1203	1205	1207	rVB	534005	650023	7.47%	1.243%
43	16.034	1262	1264	1273	rVB3	191229	568095	6.53%	1.086%
44	16.434	1297	1299	1310	rVB	278858	814846	9.37%	1.558%
45	18.046	1437	1440	1445	rVB	110829	286886	3.30%	0.548%
46	19.200	1537	1541	1550	rVB	217940	698910	8.03%	1.336%

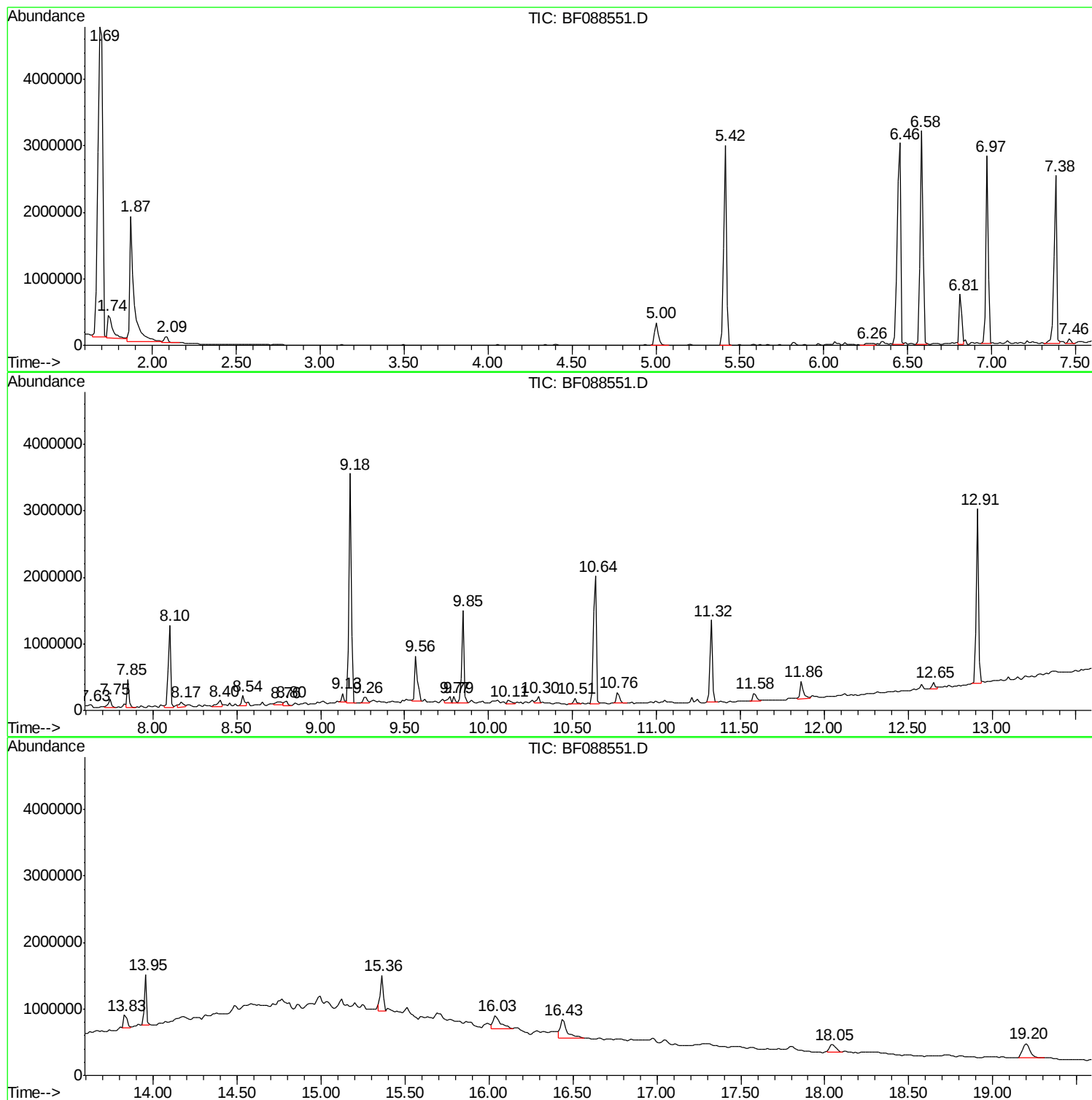
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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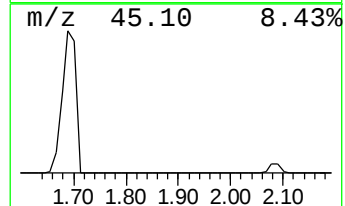
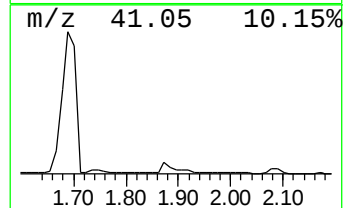
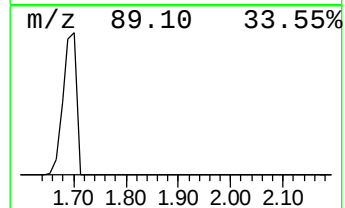
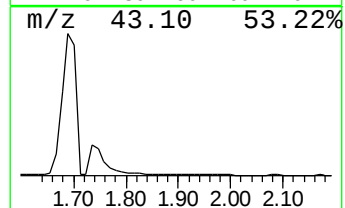
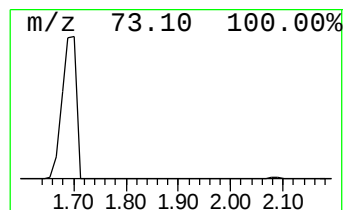
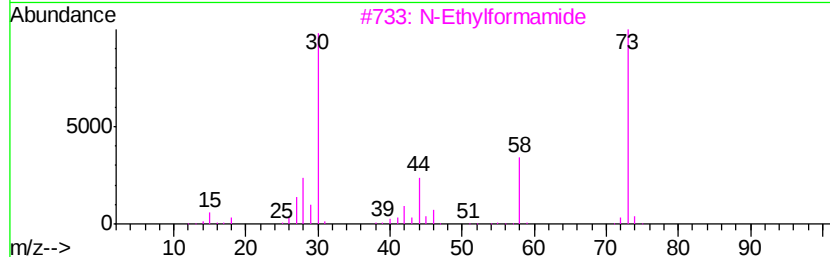
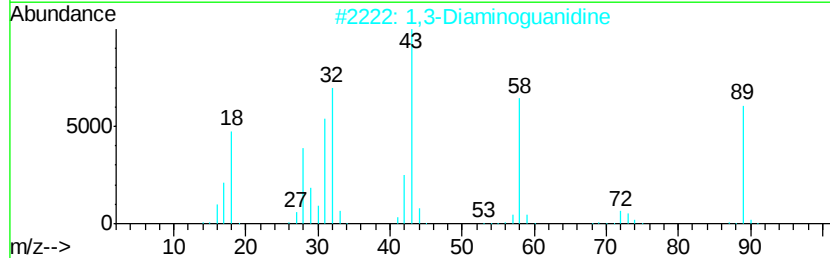
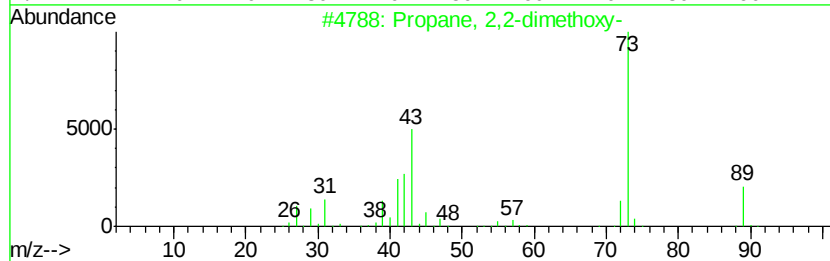
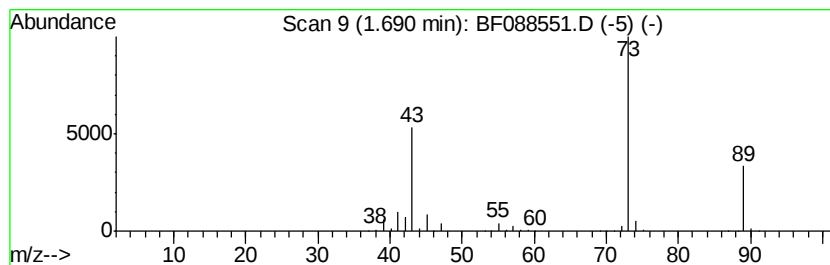
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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	202.36 ng	8699460	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	56
2		1,3-Diaminoquinidine	89	CH7N5	004364-78-7	9
3		N-Ethylformamide	73	C3H7NO	000627-45-2	9
4		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9
5		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9



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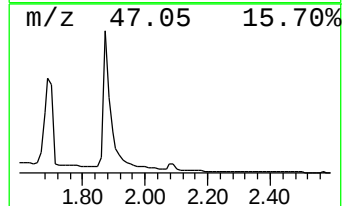
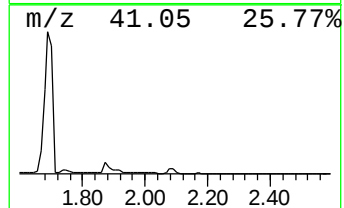
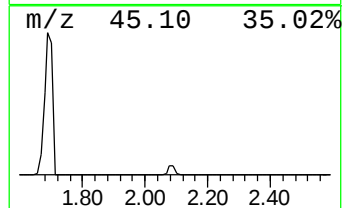
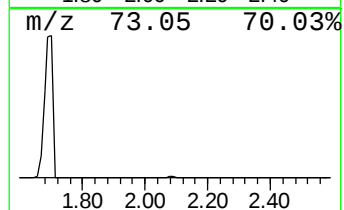
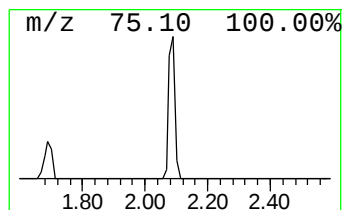
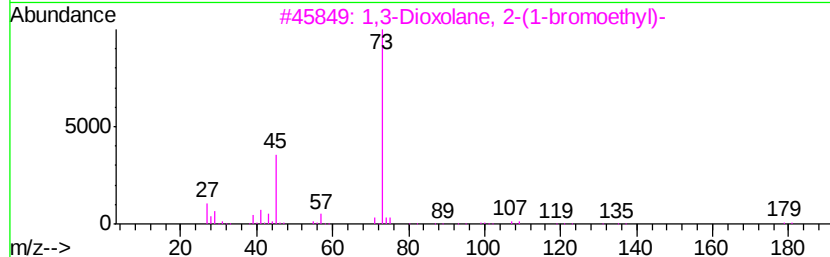
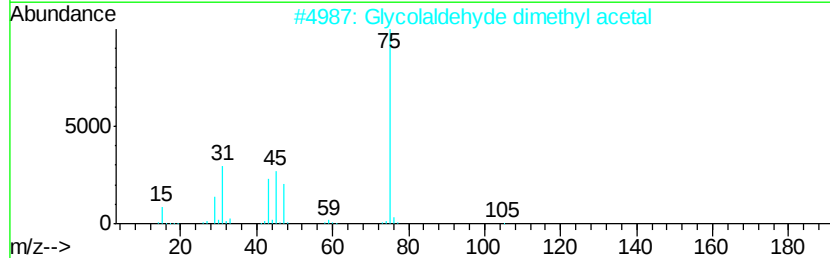
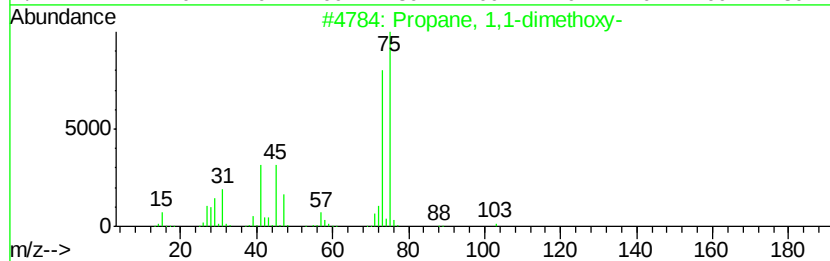
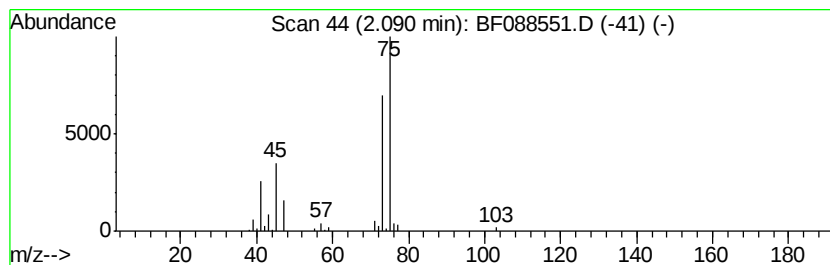
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 Propane, 1,1-dimethoxy- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.09	4.03 ng	173151	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	78
2		Glycolaldehyde dimethyl acetal	106	C4H10O3	030934-97-5	39
3		1,3-Dioxolane, 2-(1-bromoethyl)-	180	C5H9BrO2	005267-73-2	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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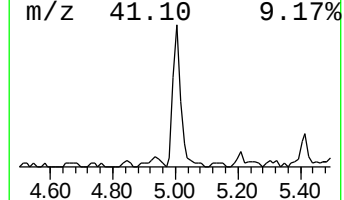
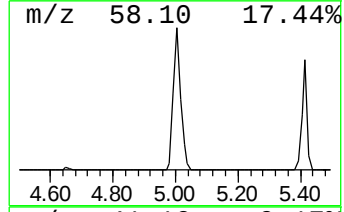
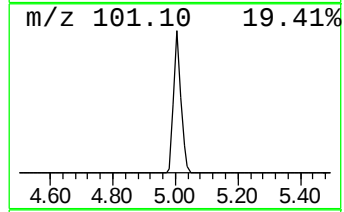
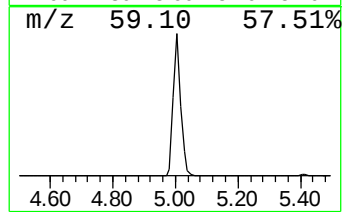
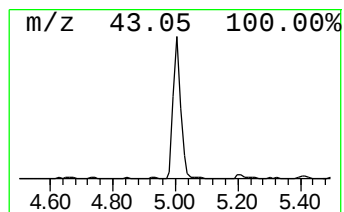
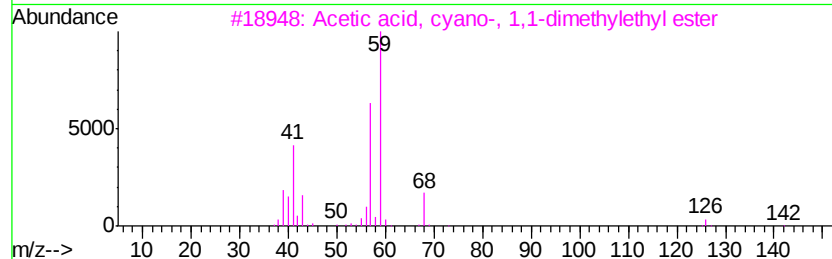
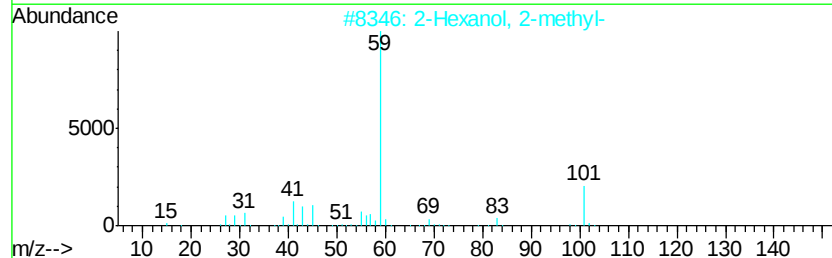
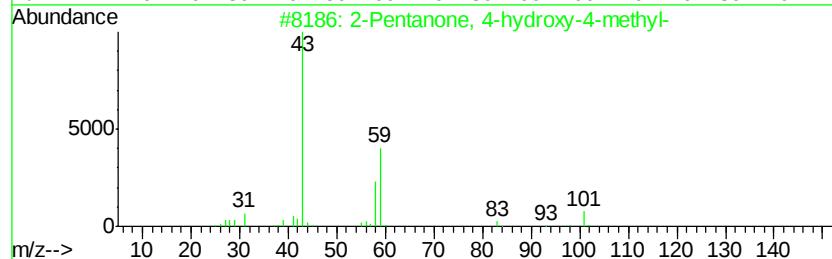
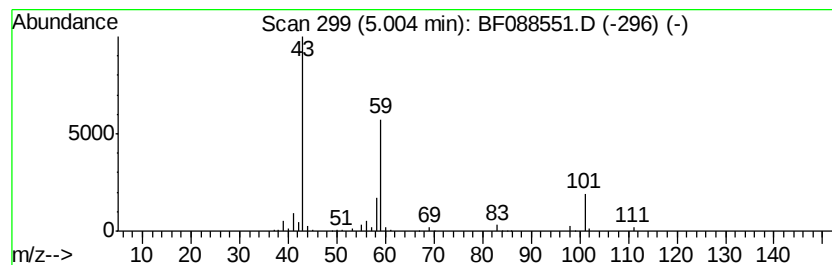
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	12.62 ng	542572	1,4-Dichlorobenzene-d4	6.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
2	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33		
3	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25		
4	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17		
5	5-Oxohexanenitrile	111	C6H9NO	010412-98-3	10		



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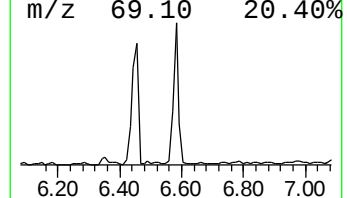
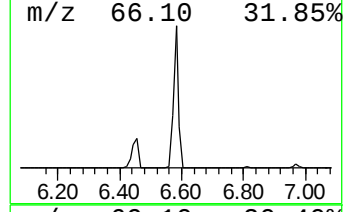
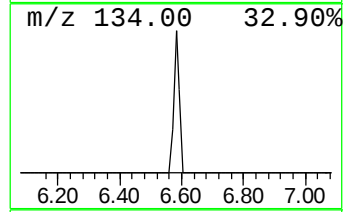
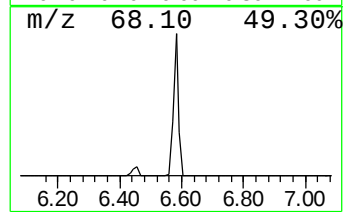
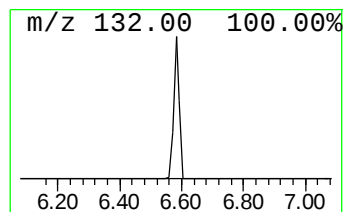
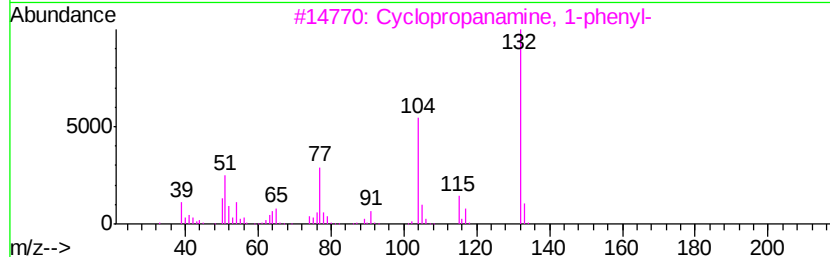
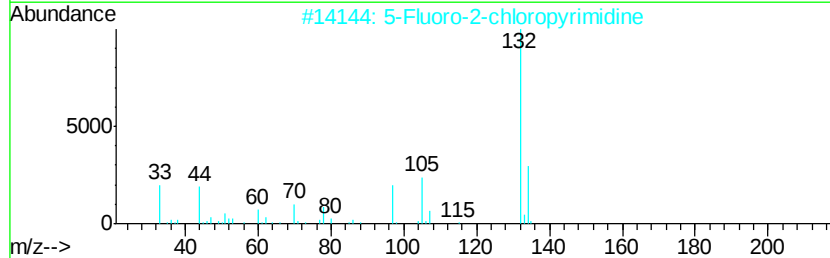
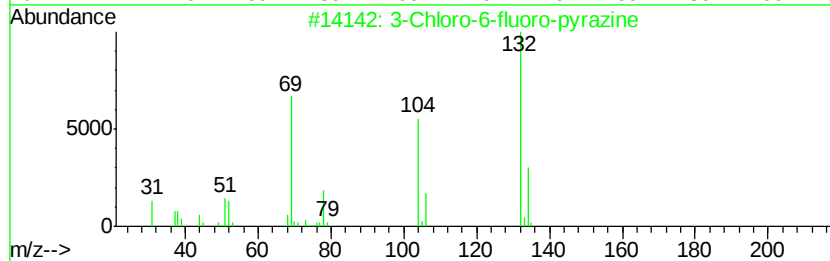
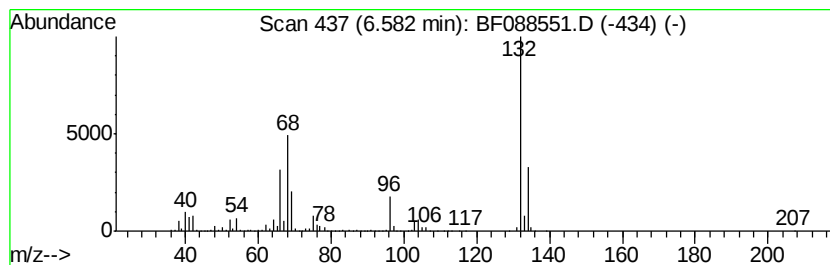
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TIC Library : C:\Database\NIST11.L
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Peak Number 6 unknown6.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	91.45 ng	3931250	1,4-Dichlorobenzene-d4	6.81

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	12
3		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10



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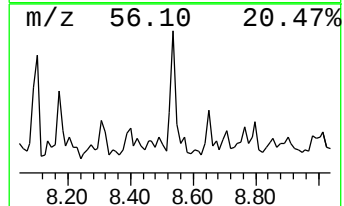
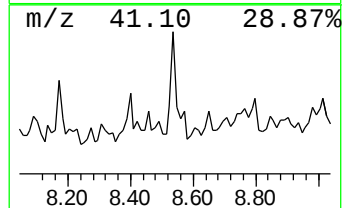
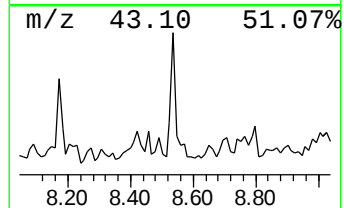
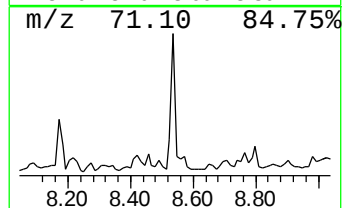
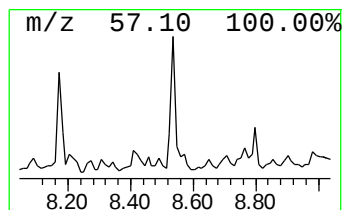
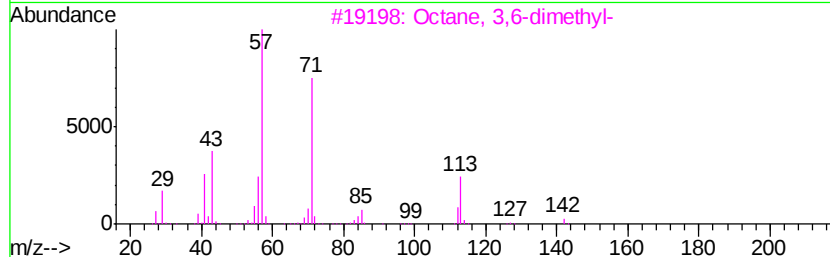
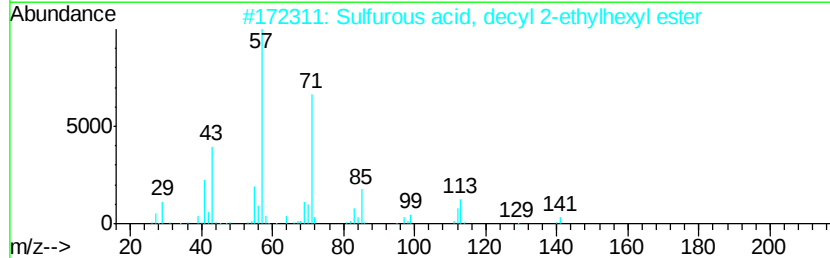
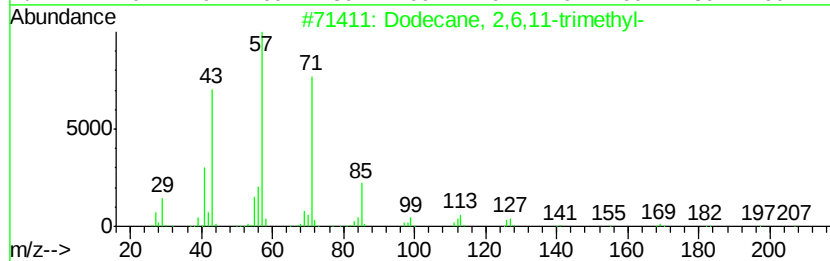
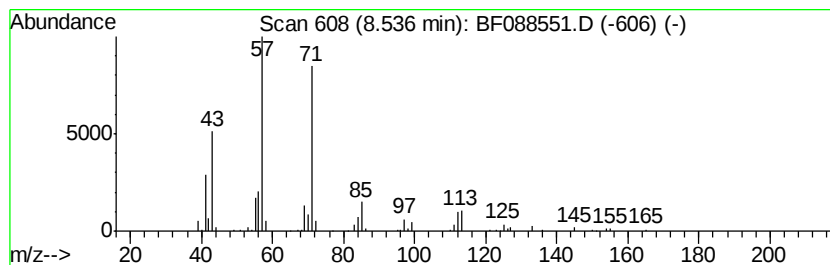
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Dodecane, 2,6,11-trimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.54	2.92 ng	189360	Napthalene-d8	8.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	72
2			Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	72
3			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	72
4			Octane, 2,3,6-trimethyl-	156	C11H24	062016-33-5	64
5			Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
Data File : BF088551.D
Acq On : 1 Jul 2016 1:48
Operator : UM/SJ
Sample : H3630-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
B1409-TP02W-1224

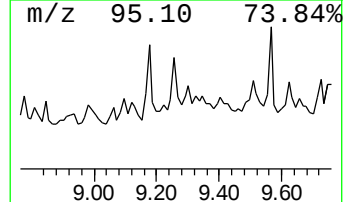
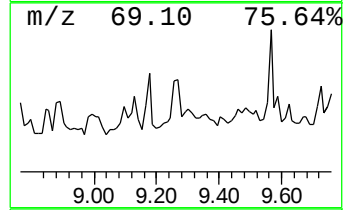
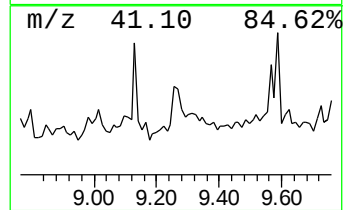
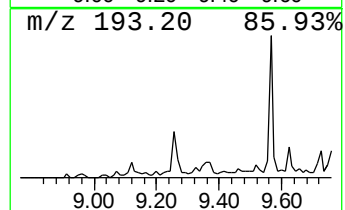
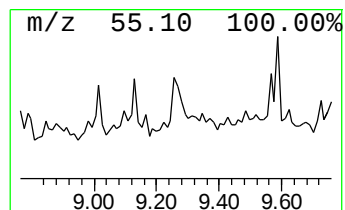
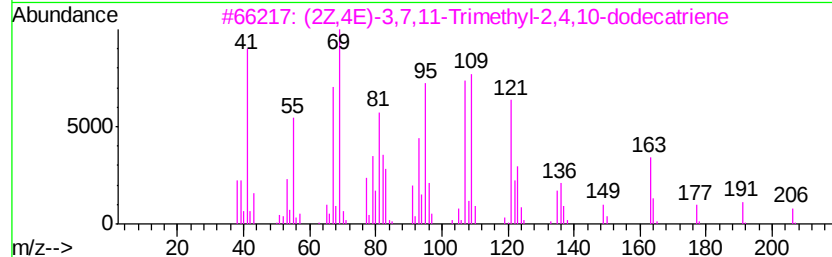
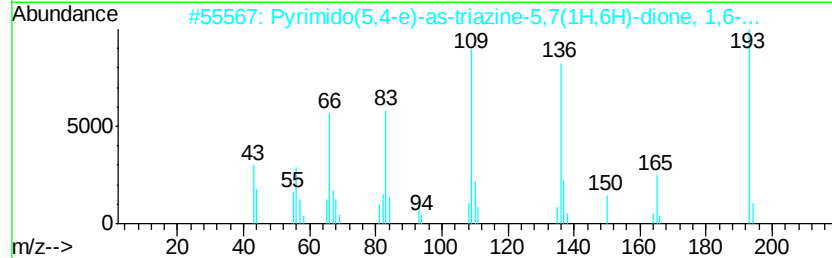
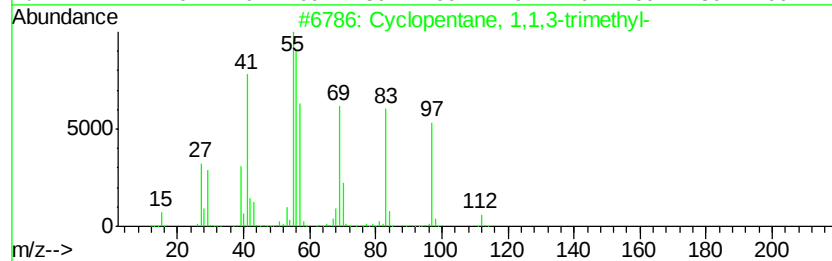
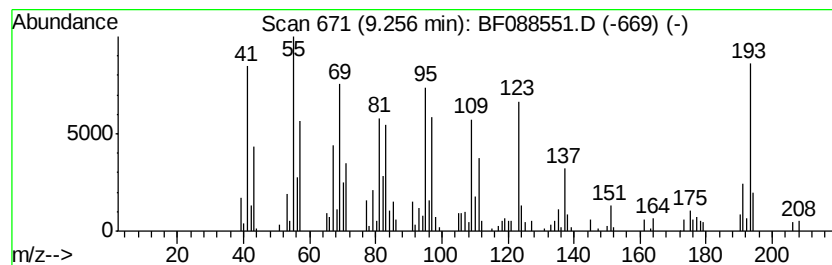
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown9.26 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.26	2.77 ng	181264	Acenaphthene-d10	9.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, 1,1,3-trimethyl-	112	C8H16	004516-69-2	45
2		Pyrimido(5,4-e)-as-triazine-5,7(...	193	C7H7N5O2	000084-82-2	35
3		(2Z,4E)-3,7,11-Trimethyl-2,4,10-...	206	C15H26	172549-29-0	25
4		3,5-Octadien-2-one, (E,E)-	124	C8H12O	030086-02-3	25
5		1-Hexyl-2-nitrocyclohexane	213	C12H23NO2	118252-04-3	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
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Operator : UM/SJ
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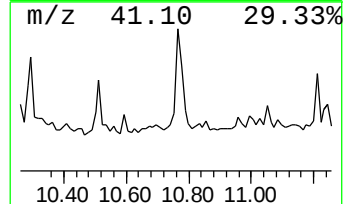
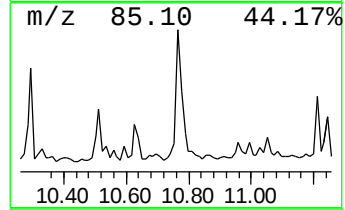
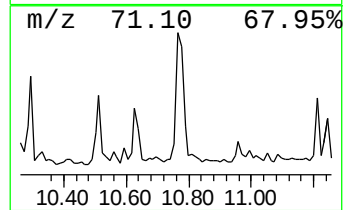
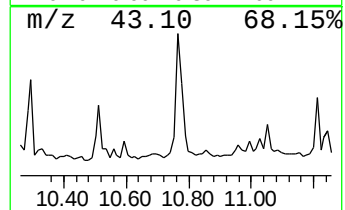
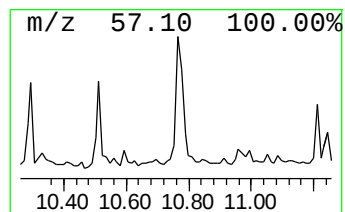
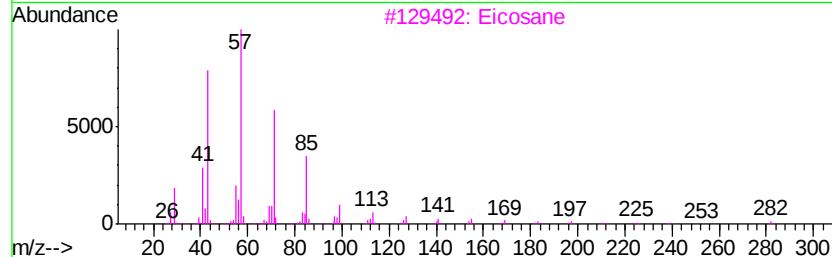
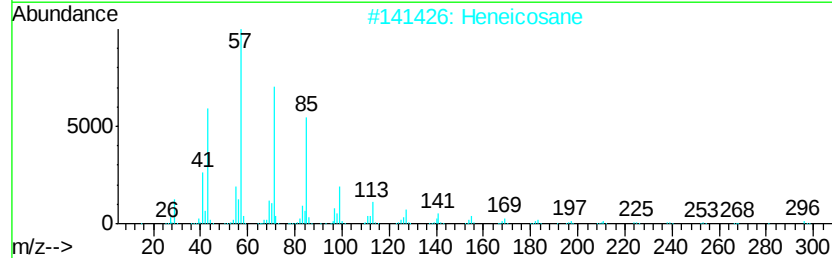
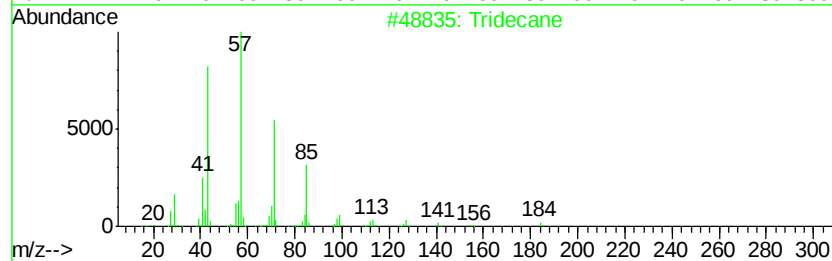
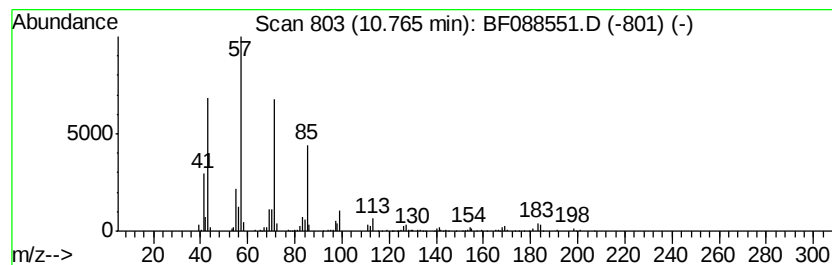
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Peak Number 12 Tridecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.76	4.57 ng	241130	Phenanthrene-d10	11.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	93
2			Heneicosane	296	C21H44	000629-94-7	87
3			Eicosane	282	C20H42	000112-95-8	87
4			Hexadecane	226	C16H34	000544-76-3	86
5			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86



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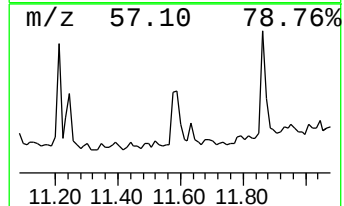
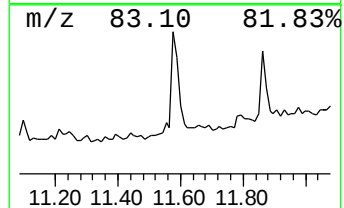
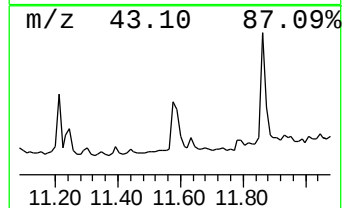
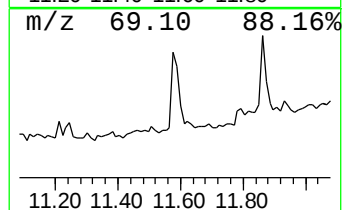
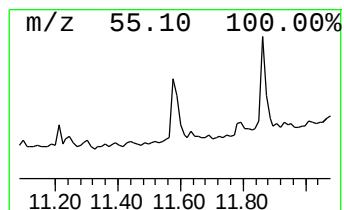
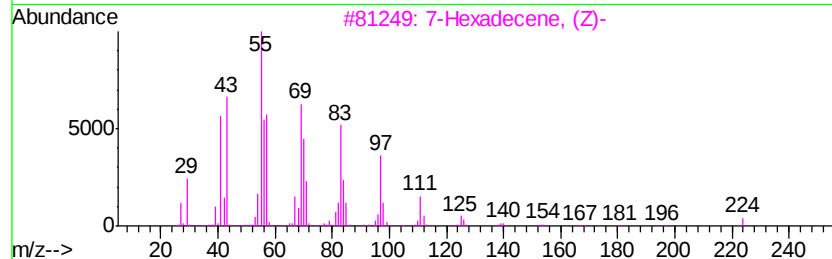
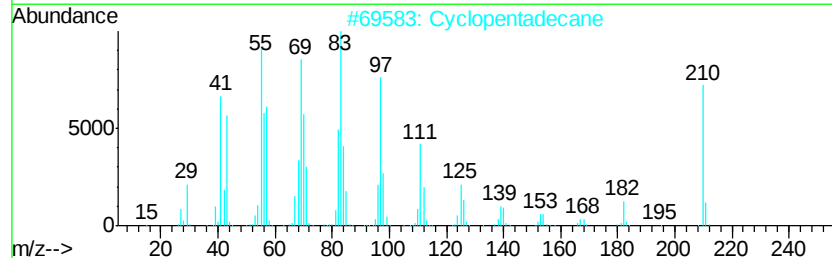
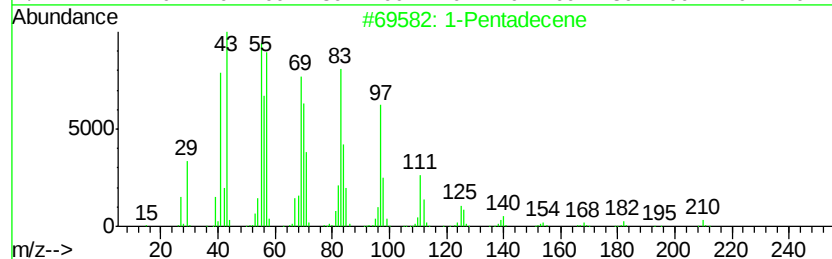
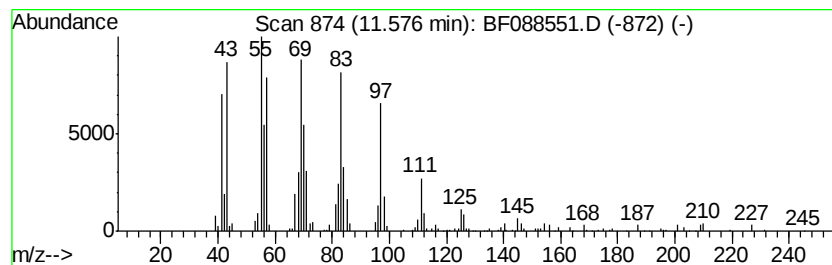
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1-Pentadecene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.58	3.36 ng	177187	Phenanthrene-d10		11.32
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentadecene	210	C15H30	013360-61-7	97
2	Cyclopentadecane	210	C15H30	000295-48-7	94
3	7-Hexadecene, (Z)-	224	C16H32	035507-09-6	91
4	n-Pentadecanol	228	C15H32O	000629-76-5	91
5	n-Heptadecanol-1	256	C17H36O	001454-85-9	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
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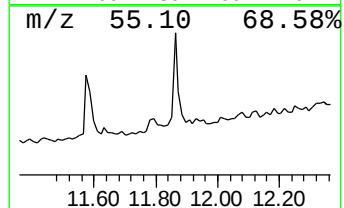
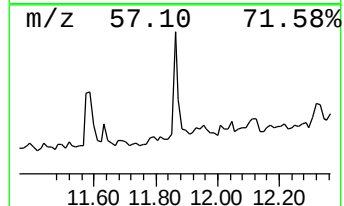
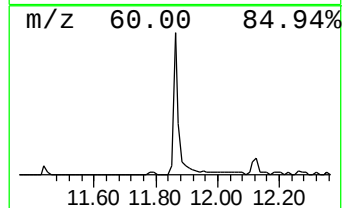
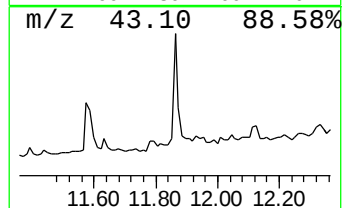
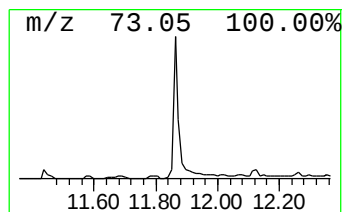
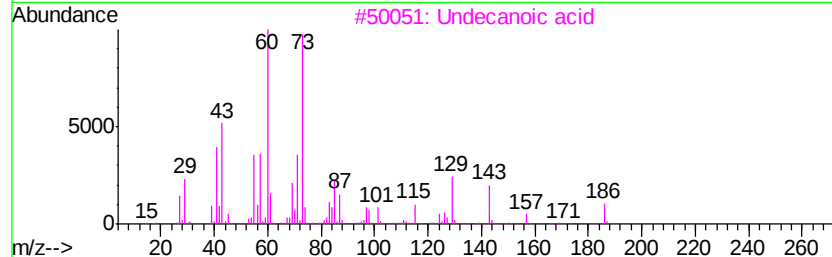
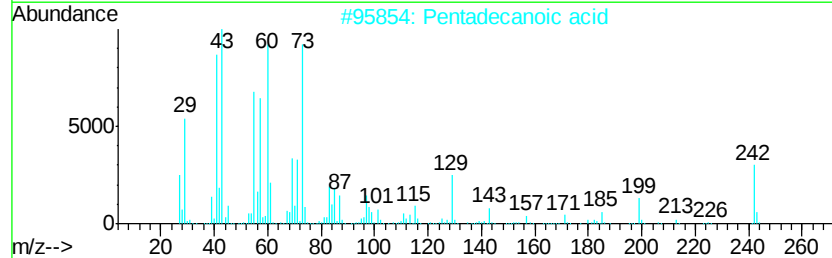
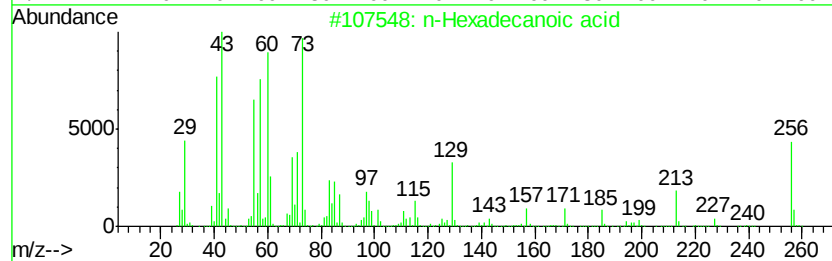
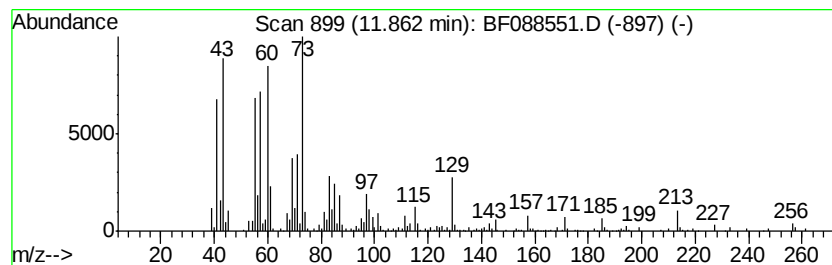
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TIC Integration Parameters: LSCINT.P

Peak Number 14 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	7.05 ng	371475	Phenanthrene-d10	11.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	83
3			Undecanoic acid	186	C11H22O2	000112-37-8	80
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5			Tridecanoic acid	214	C13H26O2	000638-53-9	58



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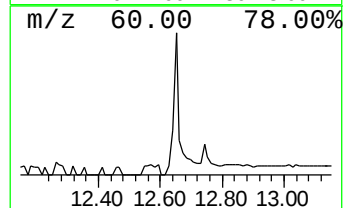
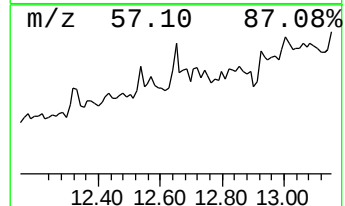
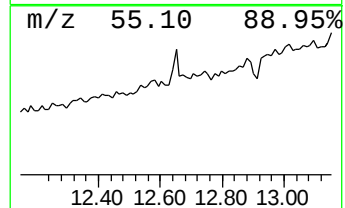
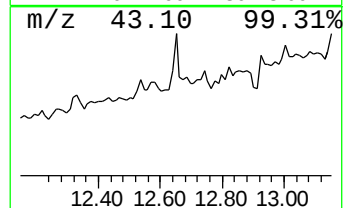
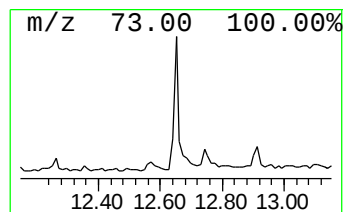
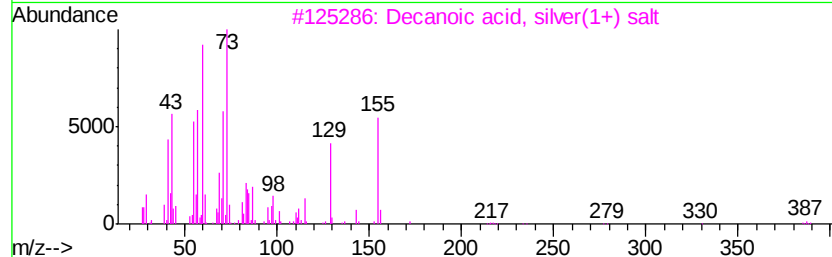
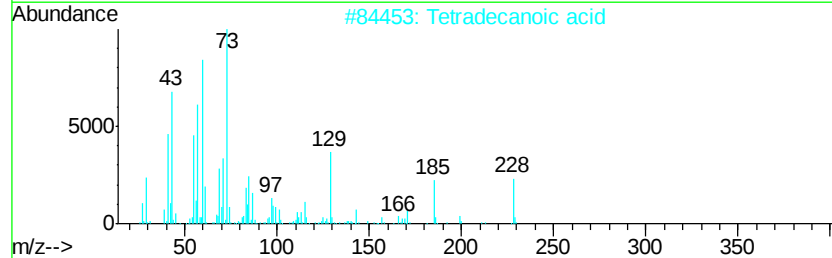
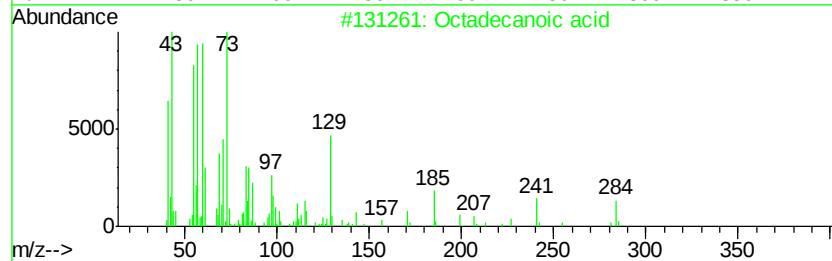
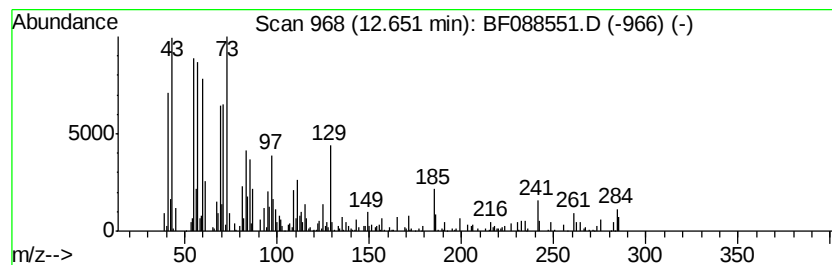
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TIC Library : C:\Database\NIST11.L
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Peak Number 15 Octadecanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	3.74 ng	129882	Chrysene-d12	13.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	98
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	47
3		Decanoic acid, silver(1+) salt	278	C10H19AgO2	013126-67-5	43
4		n-Decanoic acid	172	C10H20O2	000334-48-5	38
5		Undecanoic acid	186	C11H22O2	000112-37-8	22



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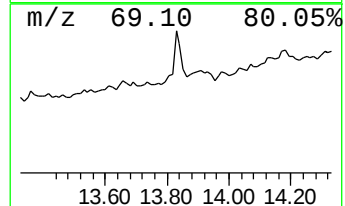
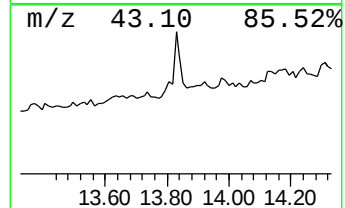
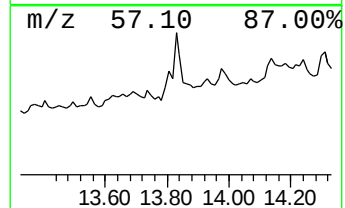
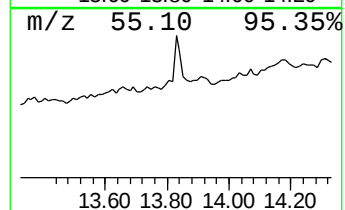
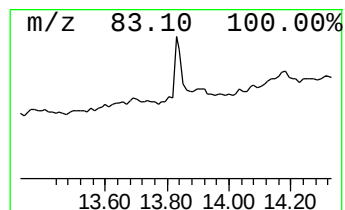
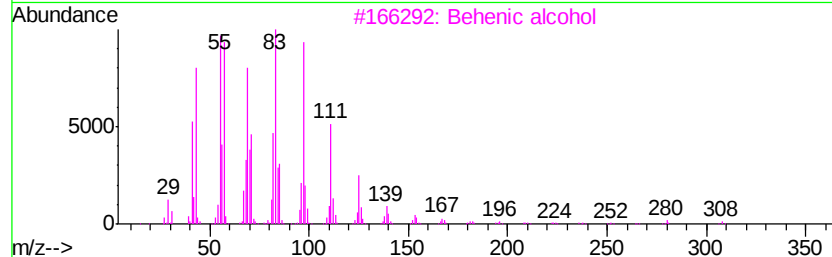
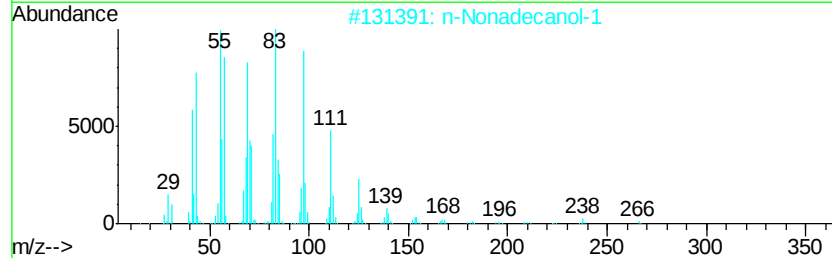
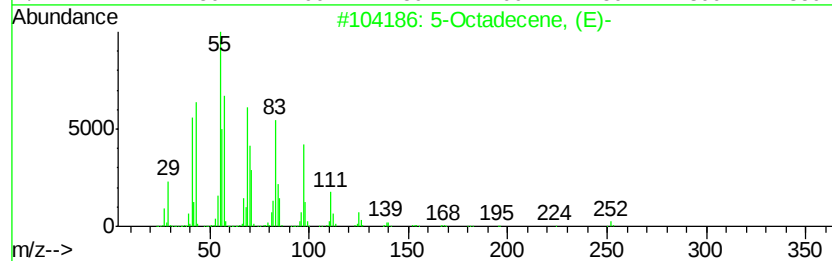
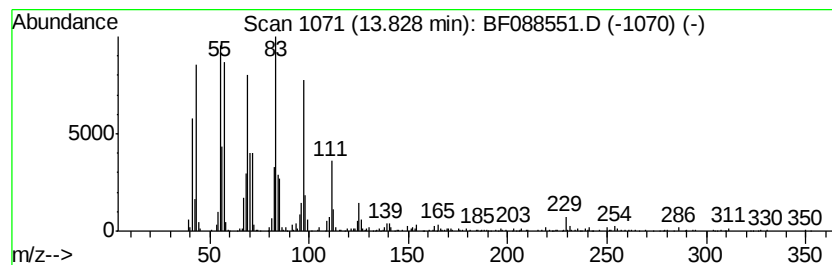
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 5-Octadecene, (E)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	6.87 ng	238642	Chrysene-d12	13.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Octadecene, (E)-	252	C18H36	007206-21-5	95
2			n-Nonadecanol-1	284	C19H40O	001454-84-8	94
3			Behenic alcohol	326	C22H46O	000661-19-8	94
4			1-Hexadecanol	242	C16H34O	036653-82-4	93
5			1-Heneicosanol	312	C21H44O	015594-90-8	91



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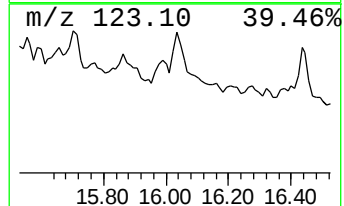
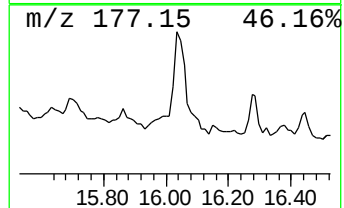
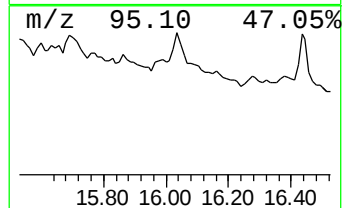
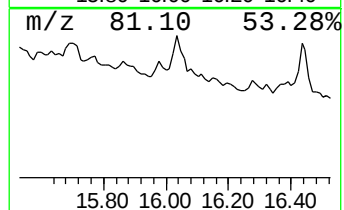
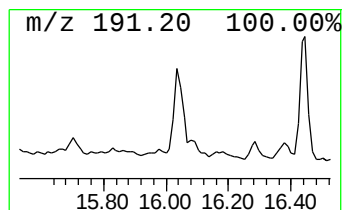
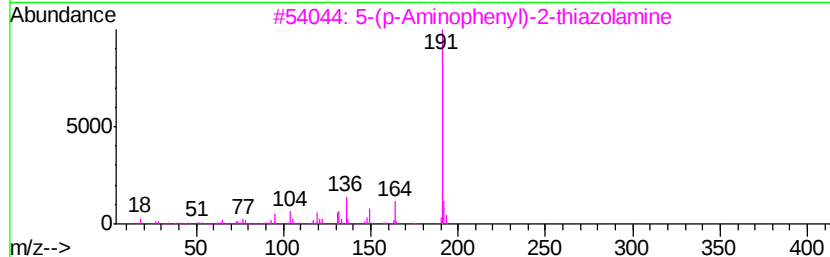
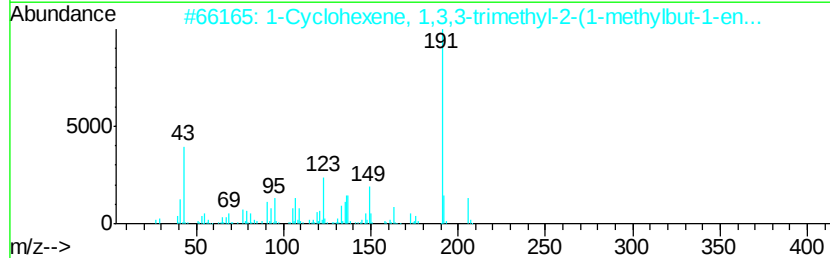
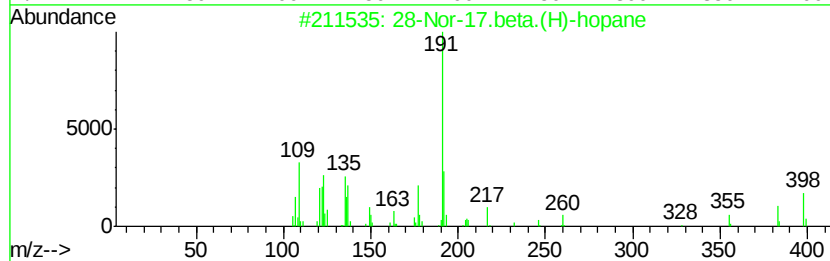
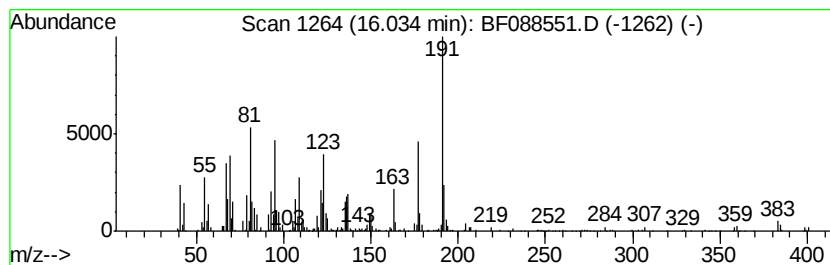
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown16.03 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.03	17.48 ng	568095	Perylene-d12	15.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	42
2		1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	35
3		5-(p-Aminophenyl)-2-thiazolamine	191	C9H9N3S	090349-87-4	25
4		(-)-Neoclovene-(II), dihydro-	206	C15H26	1000152-83-6	25
5		1-(3-Nitro-phenyl)-pyrrolidine	192	C10H12N2O2	1000300-34-7	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
Data File : BF088551.D
Acq On : 1 Jul 2016 1:48
Operator : UM/SJ
Sample : H3630-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B1409-TP02W-1224

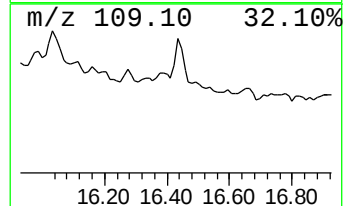
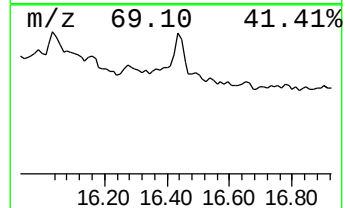
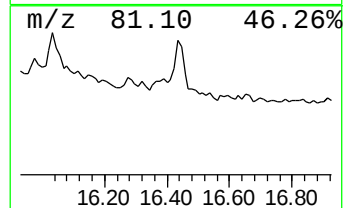
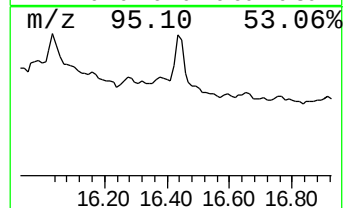
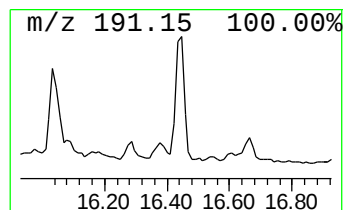
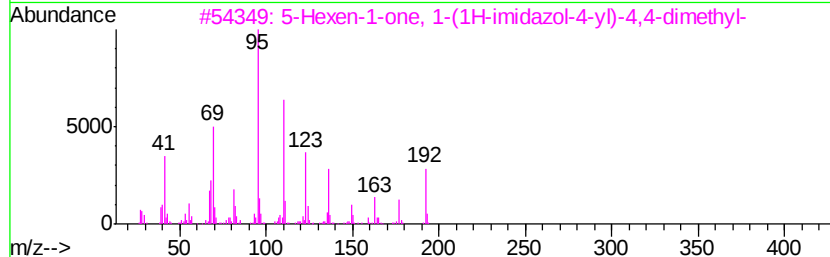
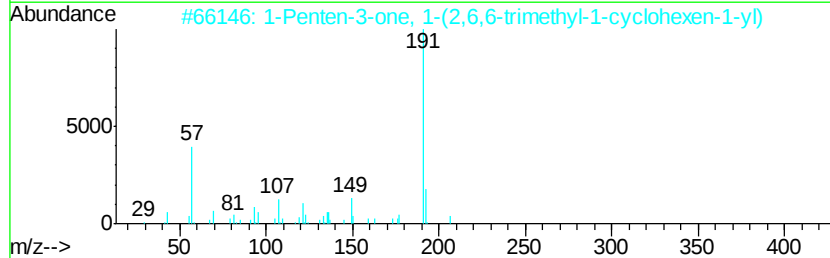
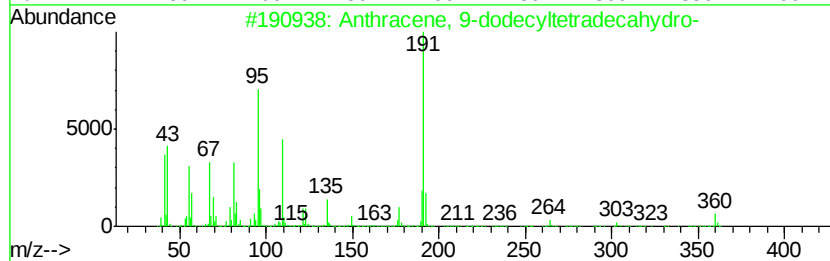
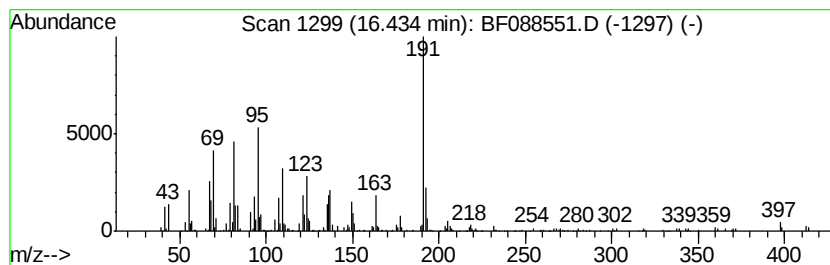
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Anthracene, 9-dodecyltetrad... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.43	25.07 ng	814846	Perylene-d12	15.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	58
2		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	43
3		5-Hexen-1-one, 1-(1H-imidazol-4-...	192	C11H16N2O	069393-41-5	42
4		Anthracene, 9-cyclohexyltetradec...	274	C20H34	055255-70-4	38
5		5-(p-Aminophenyl)-2-thiazolamine	191	C9H9N3S	090349-87-4	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
Data File : BF088551.D
Acq On : 1 Jul 2016 1:48
Operator : UM/SJ
Sample : H3630-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

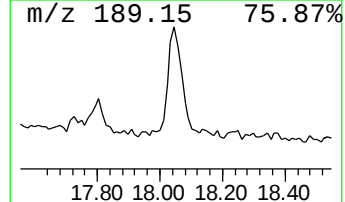
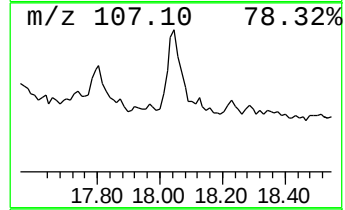
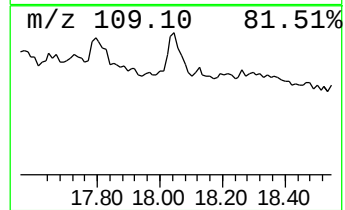
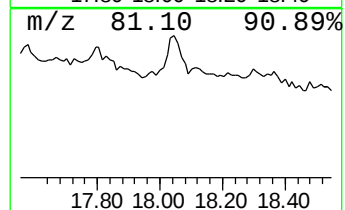
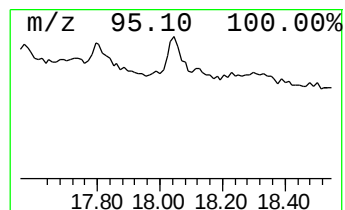
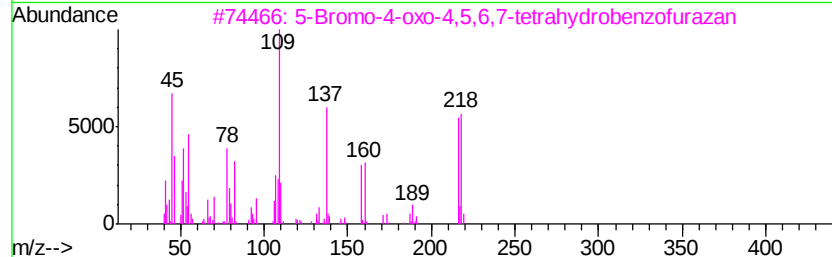
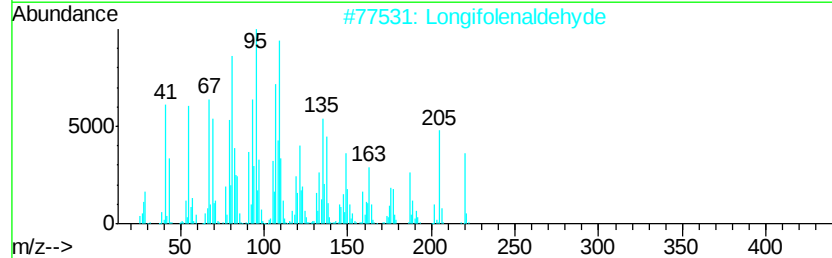
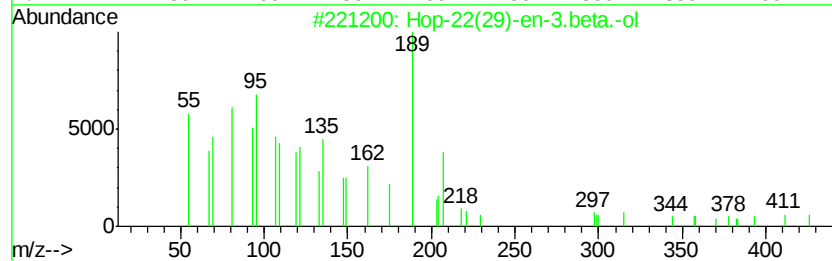
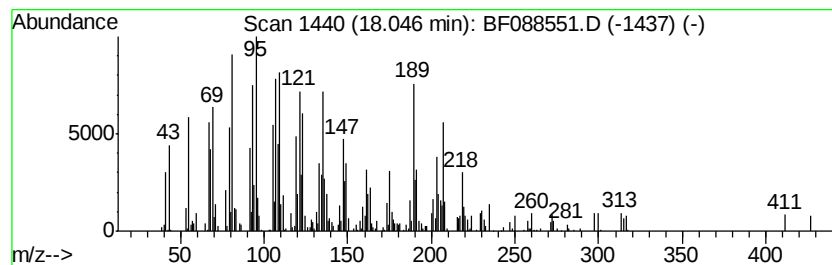
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Hop-22(29)-en-3.beta.-ol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.05	8.83 ng	286886	Perylene-d12	15.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hop-22(29)-en-3.beta.-ol	426	C30H50O	058801-23-3	78
2	Longifolenaldehyde	220	C15H24O	019890-84-7	58
3	5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	53
4	2-Methyl-3-(3-methyl-but-2-enyl)...	222	C15H26O	1000144-10-2	43
5	2-Isopropenyl-4a,8-dimethyl-1,2,...	204	C15H24	1000193-57-0	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
Data File : BF088551.D
Acq On : 1 Jul 2016 1:48
Operator : UM/SJ
Sample : H3630-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

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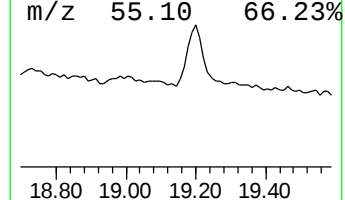
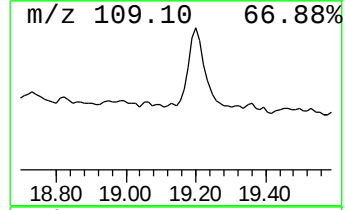
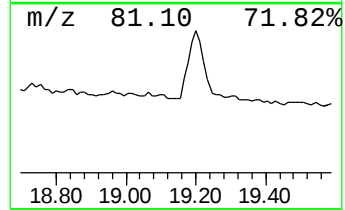
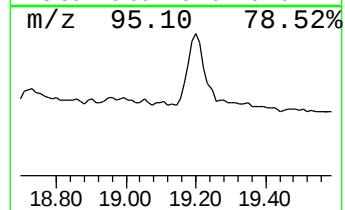
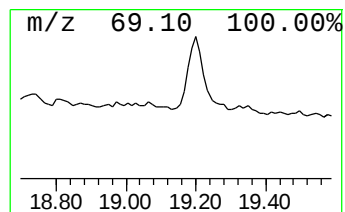
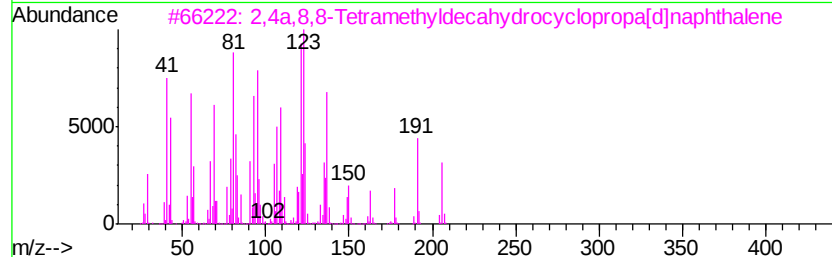
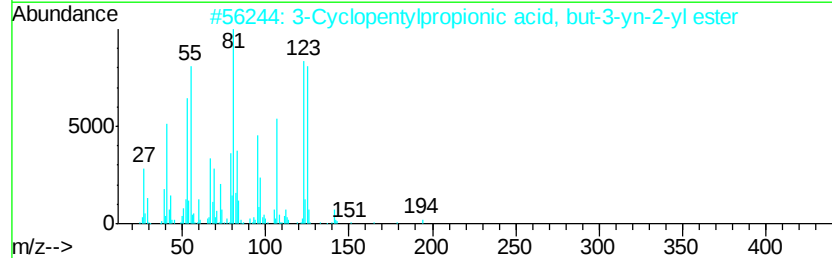
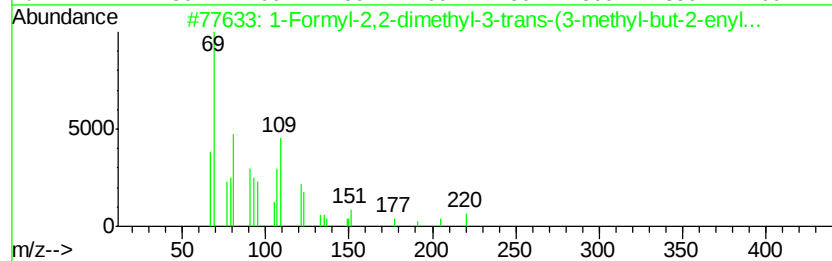
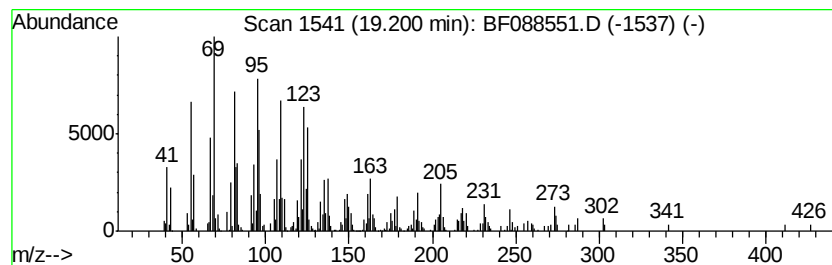
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 unknown19.20 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.20	21.50 ng	698910	Perylene-d12	15.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Formyl-2,2-dimethyl-3-trans-(3...	220	C15H24O	1000144-09-7	46
2		3-Cyclopentylpropionic acid, but...	194	C12H18O2	1000292-46-4	44
3		2,4a,8,8-Tetramethyldecahydrocyc...	206	C15H26	074022-04-1	43
4		2,3,3-Trimethyl-2-(3-methyl-buta...	206	C14H22O	1000193-60-6	42
5		Sesquirosefuran	218	C15H22O	039007-93-7	41



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063016\
Data File : BF088551.D
Acq On : 1 Jul 2016 1:48
Operator : UM/SJ
Sample : H3630-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B1409-TP02W-1224

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	202.4	ng	8699460	1	6.81	859798	20.0
Propane, 1,1-dime...	2.09	4.0	ng	173151	1	6.81	859798	20.0
2-Pentanone, 4-hy...	5.00	12.6	ng	542572	1	6.81	859798	20.0
unknown6.58	6.58	91.5	ng	3931250	1	6.81	859798	20.0
Dodecane, 2,6,11-...	8.54	2.9	ng	189360	2	8.10	1295970	20.0
unknown9.26	9.26	2.8	ng	181264	3	9.85	1307230	20.0
Tridecane	10.76	4.6	ng	241130	4	11.32	1054400	20.0
1-Pentadecene	11.58	3.4	ng	177187	4	11.32	1054400	20.0
n-Hexadecanoic acid	11.86	7.0	ng	371475	4	11.32	1054400	20.0
Octadecanoic acid	12.65	3.7	ng	129882	5	13.95	694241	20.0
5-Octadecene, (E)-	13.83	6.9	ng	238642	5	13.95	694241	20.0
unknown16.03	16.03	17.5	ng	568095	6	15.36	650023	20.0
Anthracene, 9-dod...	16.43	25.1	ng	814846	6	15.36	650023	20.0
Hop-22(29)-en-3.b...	18.05	8.8	ng	286886	6	15.36	650023	20.0
unknown19.20	19.20	21.5	ng	698910	6	15.36	650023	20.0