

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071017\
 Data File : BF096597.D
 Acq On : 10 Jul 2017 13:29
 Operator : SJ/JU
 Sample : I4086-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 A2882

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-BF070117.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.863	468	472	480	rVB	211055	250834	6.65%	0.809%
2	5.292	540	545	548	rBV	1706127	1679987	44.56%	5.418%
3	5.504	577	581	586	rBV2	57371	68097	1.81%	0.220%
4	6.316	715	719	726	rVB	1312663	1140835	30.26%	3.679%
5	6.398	730	733	737	rVB	43224	41136	1.09%	0.133%
6	6.451	737	742	745	rBV	2835025	2795319	74.14%	9.015%
7	6.681	777	781	785	rBV	894121	766431	20.33%	2.472%
8	6.751	789	793	799	rBV2	90078	102691	2.72%	0.331%
9	6.839	799	808	811	rBV	2510955	2501736	66.35%	8.069%
10	7.086	845	850	856	rVB	40768	43724	1.16%	0.141%
11	7.245	871	877	880	rBV	2189553	2330136	61.80%	7.515%
12	7.281	880	883	890	rVV2	41665	49529	1.31%	0.160%
13	7.351	890	895	901	rVB5	25697	46335	1.23%	0.149%
14	7.498	915	920	924	rVB2	36237	46583	1.24%	0.150%
15	7.710	952	956	959	rVB2	44959	41637	1.10%	0.134%
16	7.875	977	984	987	rBV	58810	69177	1.83%	0.223%
17	7.963	995	999	1002	rBV	1145272	1080564	28.66%	3.485%
18	7.992	1002	1004	1009	rVV	244041	203664	5.40%	0.657%
19	8.039	1009	1012	1015	rVB	74288	69988	1.86%	0.226%
20	8.216	1036	1042	1044	rBV	84558	113511	3.01%	0.366%
21	8.380	1067	1070	1075	rVB2	44359	59180	1.57%	0.191%
22	8.669	1114	1119	1123	rBV3	58668	67971	1.80%	0.219%
23	9.045	1177	1183	1186	rBV	3419952	3770443	100.00%	12.160%
24	9.169	1201	1204	1207	rBV3	33627	39738	1.05%	0.128%
25	9.374	1234	1239	1242	rBV5	29874	43018	1.14%	0.139%
26	9.433	1245	1249	1254	rVB2	259559	250589	6.65%	0.808%
27	9.716	1291	1297	1300	rBV	1243586	1203028	31.91%	3.880%
28	9.845	1316	1319	1321	rBV	48175	47159	1.25%	0.152%
29	9.869	1321	1323	1325	rVV	56552	46056	1.22%	0.149%
30	9.927	1328	1333	1337	rVV4	62310	91808	2.43%	0.296%
31	9.980	1340	1342	1347	rVB2	36196	39756	1.05%	0.128%
32	10.280	1391	1393	1397	rVB	49233	40284	1.07%	0.130%
33	10.504	1425	1431	1434	rBV	2019048	2208479	58.57%	7.123%
34	10.586	1441	1445	1450	rBV	105327	103396	2.74%	0.333%

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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	10.633	1450	1453	1465	rVB	43611	54851	1.45%	0.177%
36	11.051	1520	1524	1526	rBV	47979	48924	1.30%	0.158%
37	11.127	1532	1537	1540	rBV2	37027	67716	1.80%	0.218%
38	11.192	1543	1548	1551	rVV	1294770	1128329	29.93%	3.639%
39	11.480	1594	1597	1600	rVB2	56366	49419	1.31%	0.159%
40	11.733	1631	1640	1643	rBV	162403	165824	4.40%	0.535%
41	11.763	1643	1645	1648	rVB	143867	107448	2.85%	0.347%
42	12.145	1704	1710	1712	rBV	43440	45611	1.21%	0.147%
43	12.168	1712	1714	1720	rVB2	45425	43369	1.15%	0.140%
44	12.245	1720	1727	1730	rBV5	17543	41323	1.10%	0.133%
45	12.474	1763	1766	1769	rVB	173478	142806	3.79%	0.461%
46	12.786	1810	1819	1822	rBV	2810663	3558449	94.38%	11.477%
47	13.121	1872	1876	1877	rBV	44847	44330	1.18%	0.143%
48	13.145	1877	1880	1885	rVB	70648	76127	2.02%	0.246%
49	13.674	1965	1970	1975	rBV3	463943	740869	19.65%	2.389%
50	13.821	1991	1995	1998	rBV	1182508	1074620	28.50%	3.466%
51	14.004	2021	2026	2031	rVB3	111851	161721	4.29%	0.522%
52	14.474	2100	2106	2112	rBV	435988	605966	16.07%	1.954%
53	14.792	2154	2160	2166	rVB	80301	121101	3.21%	0.391%
54	15.215	2226	2232	2237	rBV	666263	760067	20.16%	2.451%
55	15.268	2237	2241	2244	rVV2	23861	45605	1.21%	0.147%
56	15.321	2244	2250	2258	rVB	175657	321675	8.53%	1.037%
57	15.756	2321	2324	2329	rVB2	24581	39610	1.05%	0.128%
58	16.527	2448	2455	2461	rVB	40820	107241	2.84%	0.346%

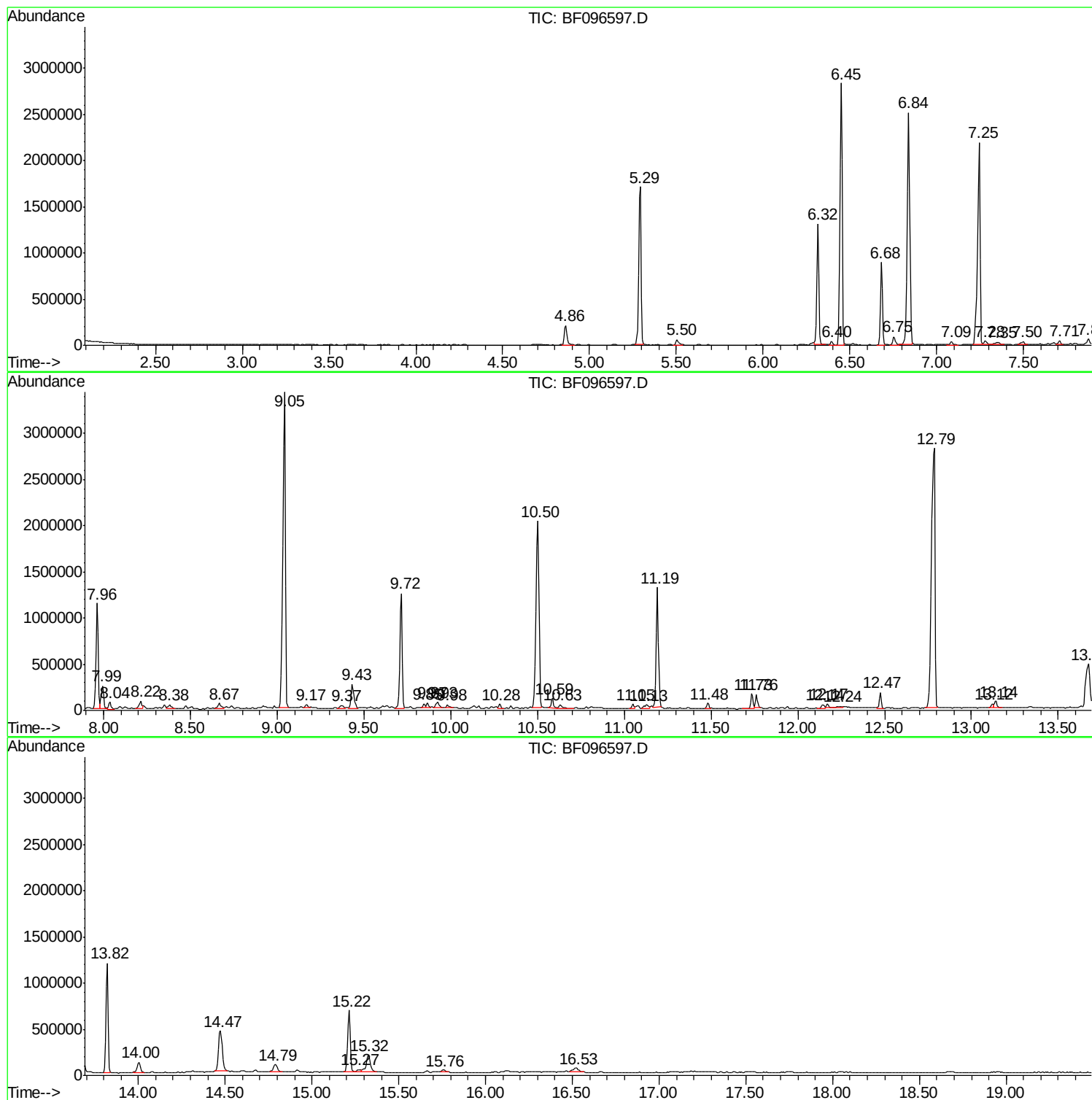
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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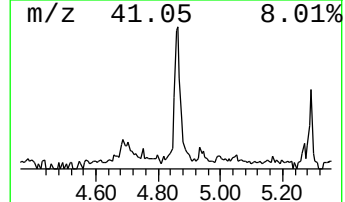
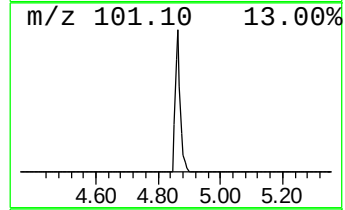
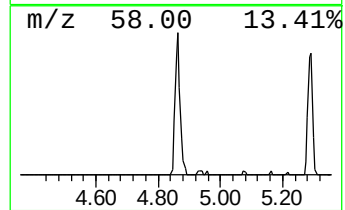
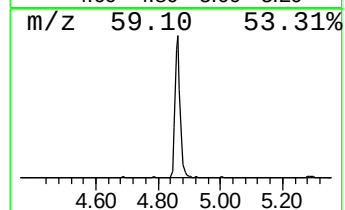
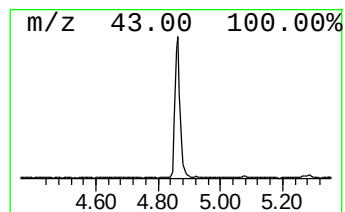
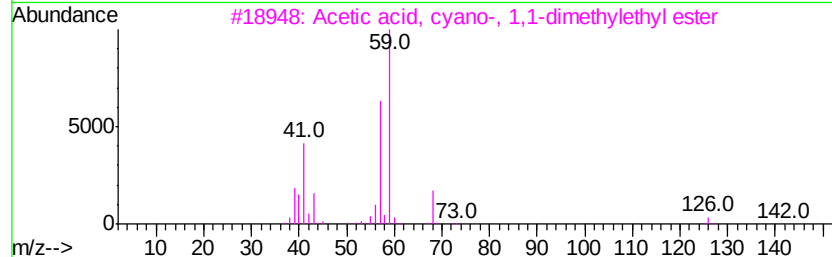
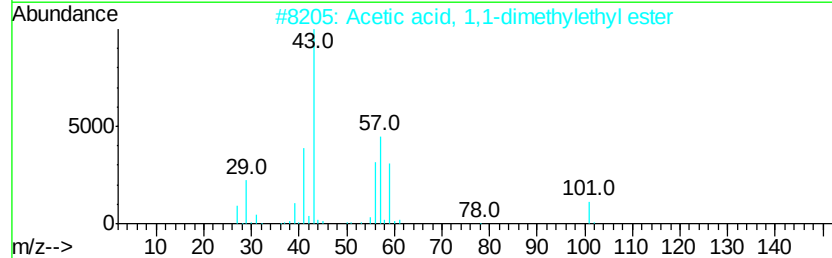
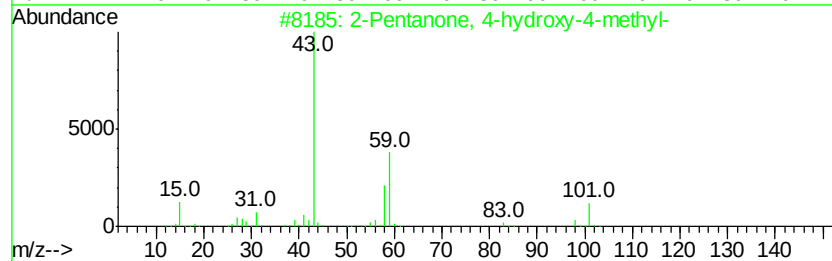
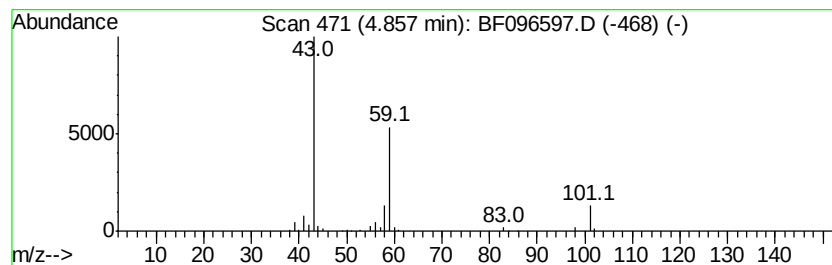
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF070117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	6.55 ng	250834	1,4-Dichlorobenzene-d4	6.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	37
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33



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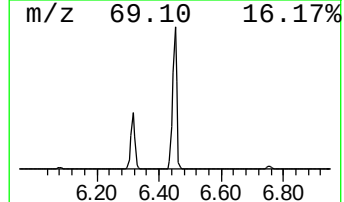
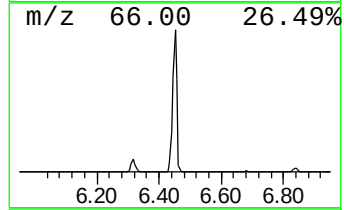
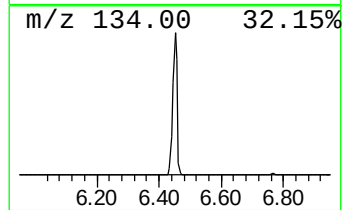
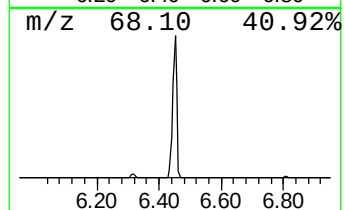
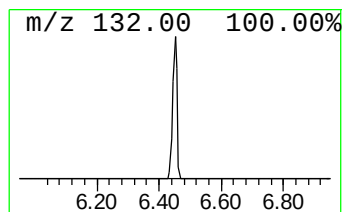
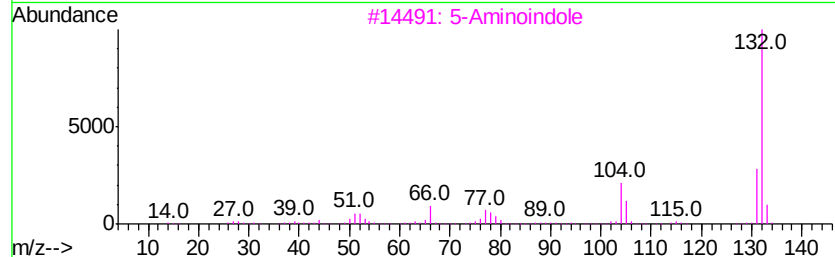
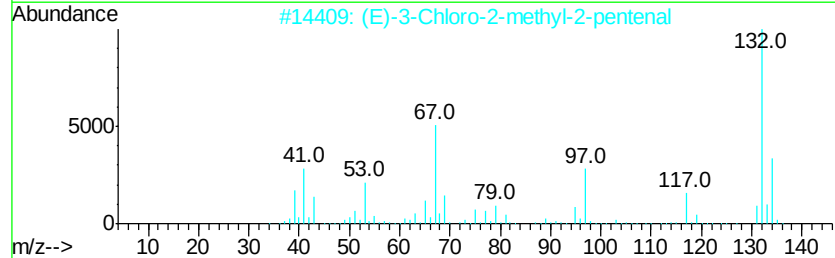
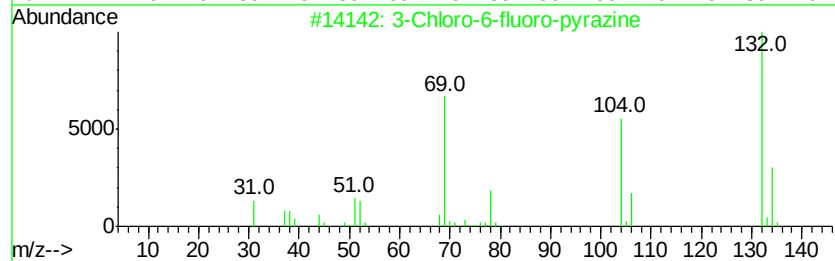
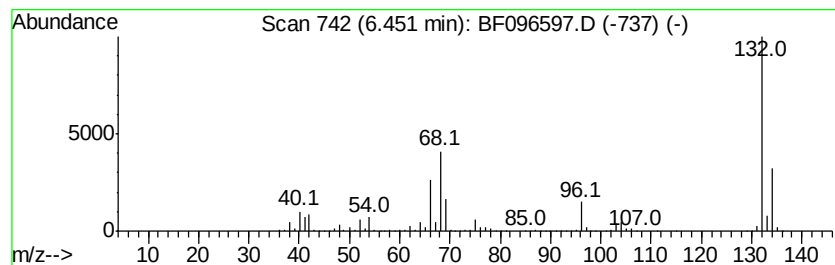
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.45 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.45	72.94 ng	2795320	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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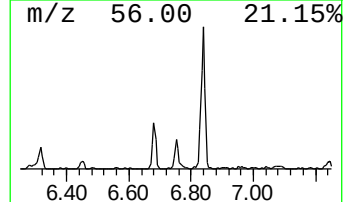
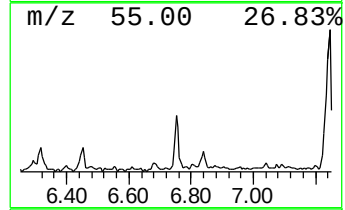
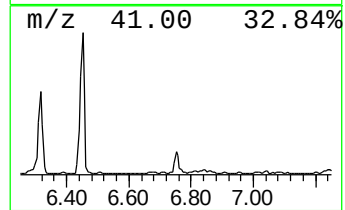
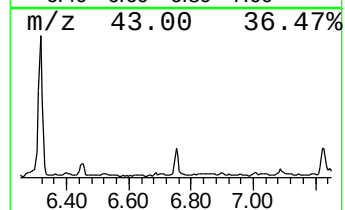
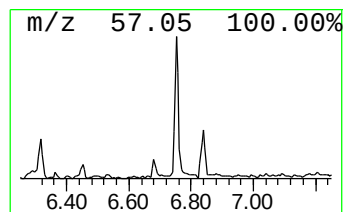
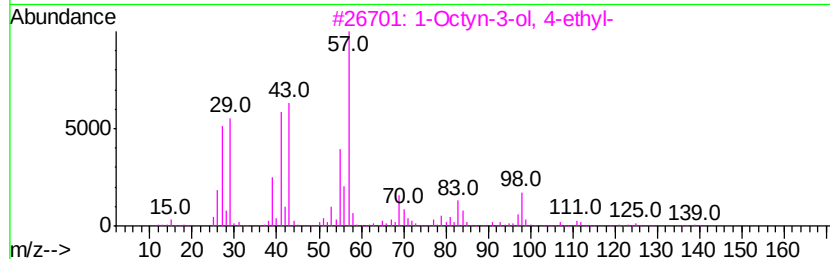
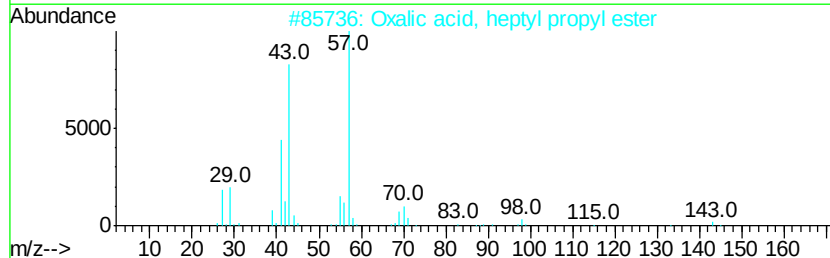
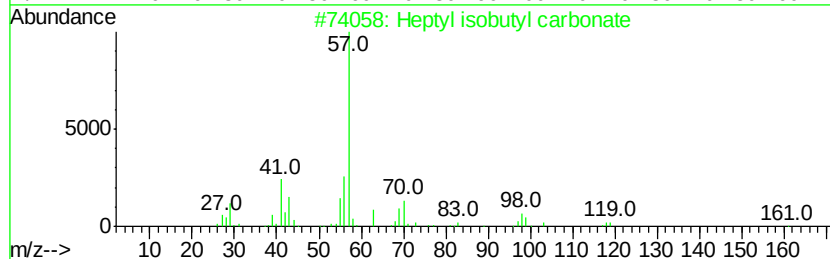
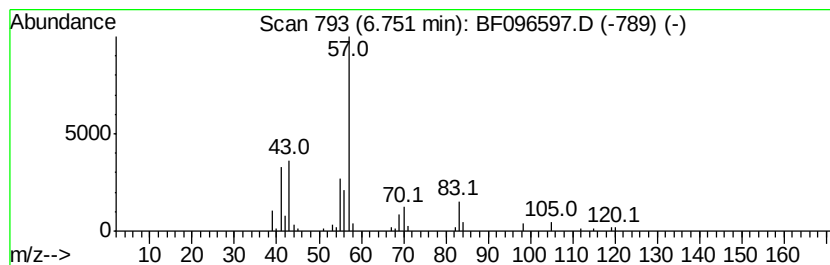
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Heptyl isobutyl carbonate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	2.68 ng	102691	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptyl isobutyl carbonate	216	C12H24O3	959068-08-7	53
2		Oxalic acid, heptyl propyl ester	230	C12H22O4	1000309-26-0	43
3		1-Octyn-3-ol, 4-ethyl-	154	C10H18O	005877-42-9	42
4		1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	42
5		Heptane, 2,3,6-trimethyl-	142	C10H22	004032-93-3	42



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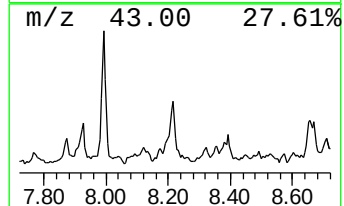
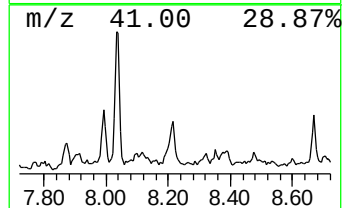
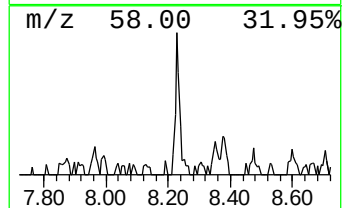
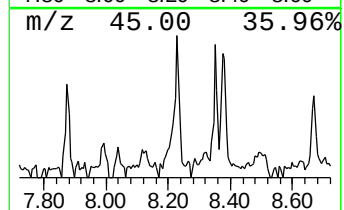
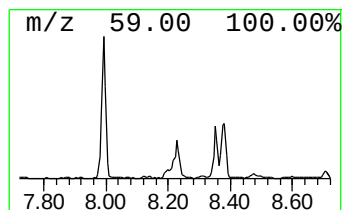
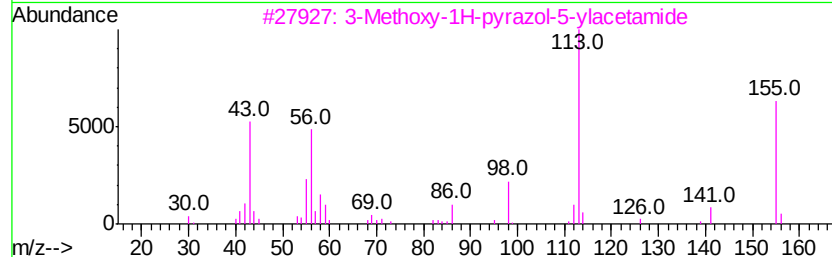
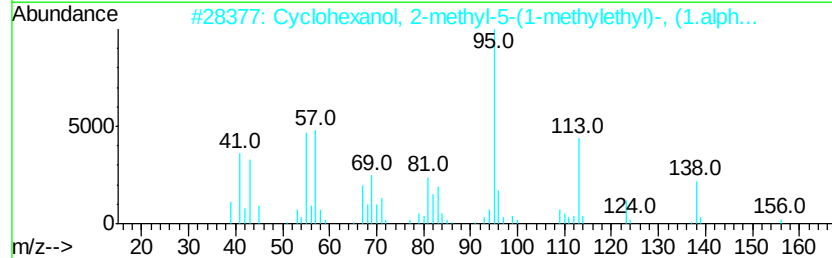
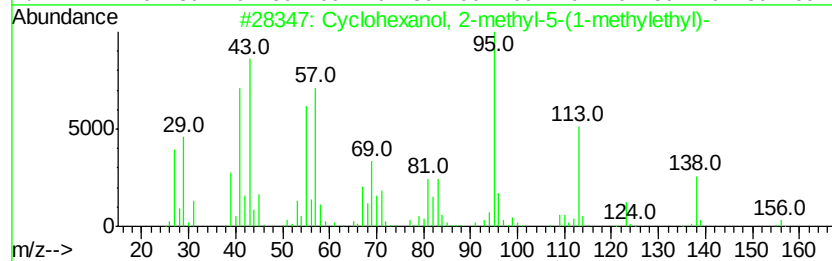
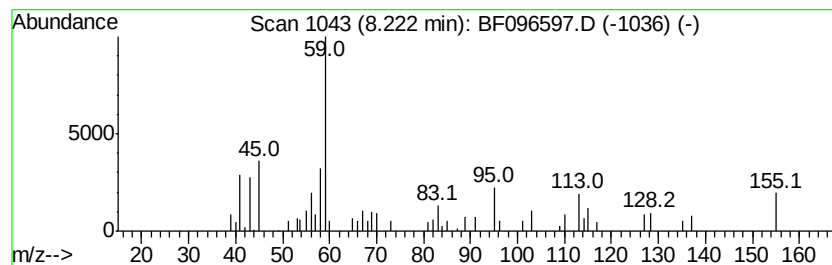
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 unknown8.22 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.22	2.10 ng	113511	Naphthalene-d8	7.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanol, 2-methyl-5-(1-meth...	156	C10H20O	060320-28-7	27
2		Cyclohexanol, 2-methyl-5-(1-meth...	156	C10H20O	000499-69-4	27
3		3-Methoxy-1H-pyrazol-5-ylacetamide	155	C6H9N3O2	1000284-55-3	18
4		2,3-Dimethyl-isoxazol-5(2H)-one	113	C5H7N02	007713-67-9	14
5		Cyclohexanecarboxylic acid, 2-(i...	282	C9H15IO2	120790-29-6	14



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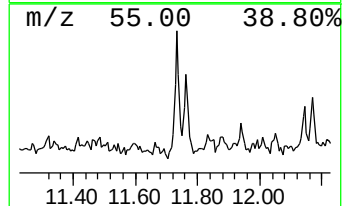
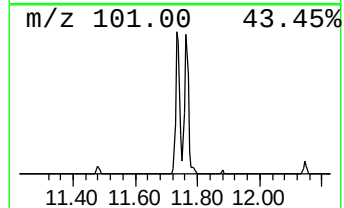
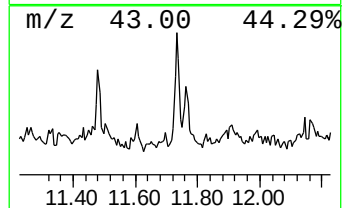
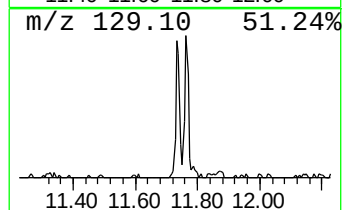
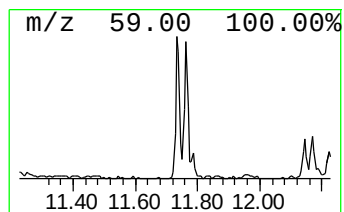
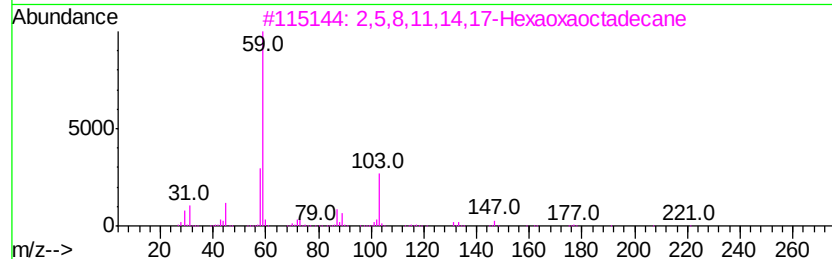
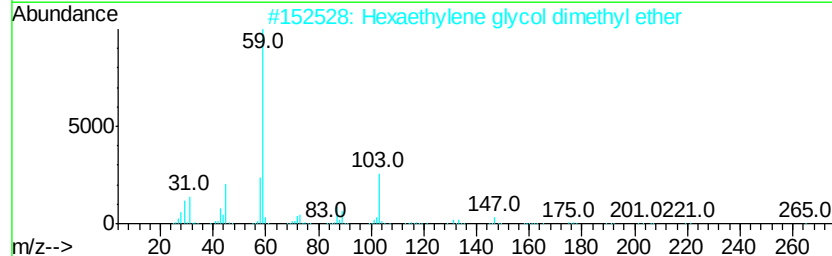
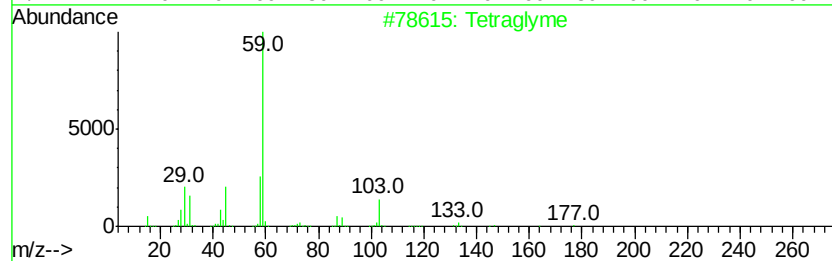
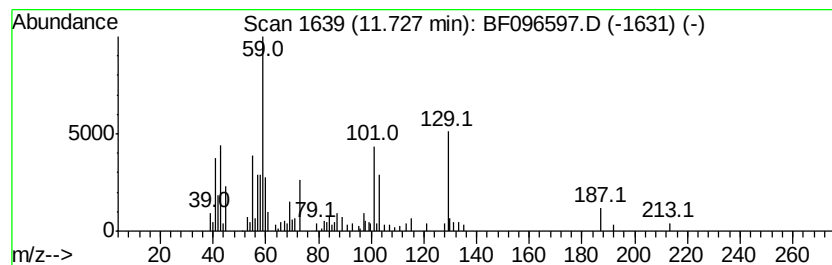
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown11.73 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.73	2.94 ng	165824	Phenanthrene-d10	11.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetraolyme	222	C10H22O5	000143-24-8	43
2		Hexaethylene glycol dimethyl ether	310	C14H30O7	001072-40-8	43
3		2,5,8,11,14,17-Hexaoxaoctadecane	266	C12H26O6	001191-87-3	43
4		Butyric acid, 3-bromo-, methyl e...	180	C5H9BrO2	021249-59-2	38
5		2,2,3,3,5,8,11-Heptamethyl-4,7,1...	348	C18H40O4Si	1000367-07-2	37



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071017\
Data File : BF096597.D
Acq On : 10 Jul 2017 13:29
Operator : SJ/JU
Sample : I4086-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

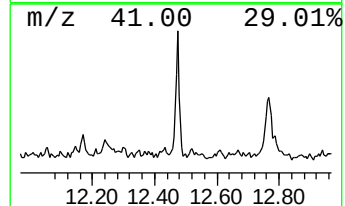
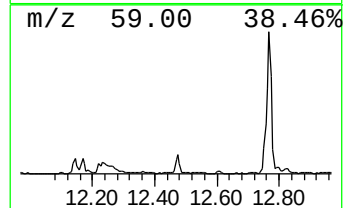
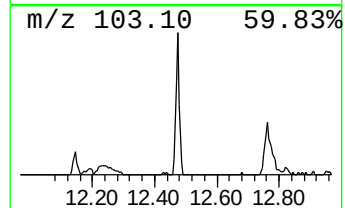
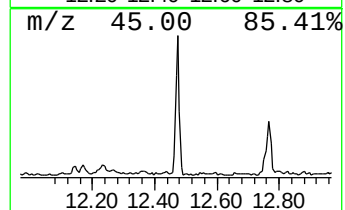
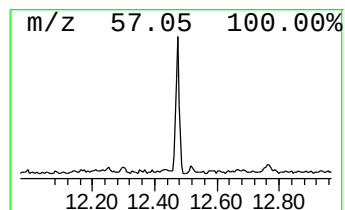
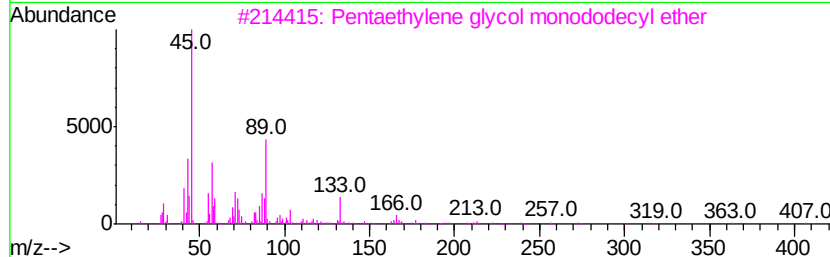
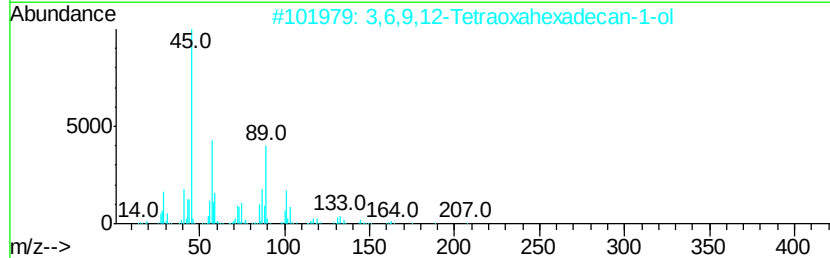
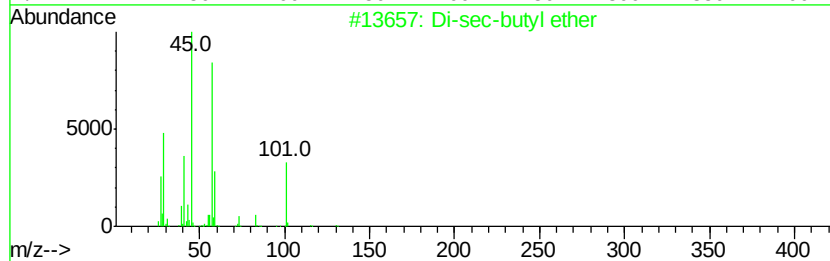
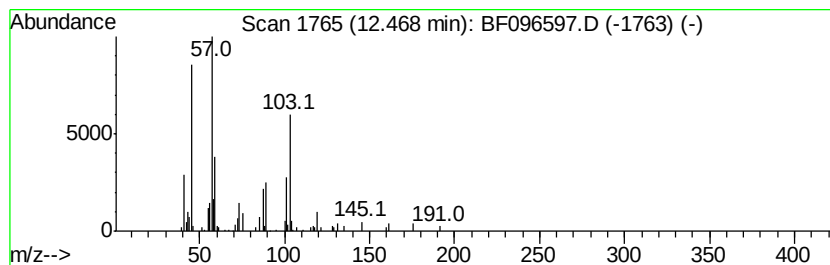
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown12.47 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.47	2.53 ng	142806	Phenanthrene-d10	11.19	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Di-sec-butyl ether	130	C8H18O	006863-58-7	47
2	3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	43
3	Pentaethylene glycol monododecyl...	406	C22H46O6	003055-95-6	38
4	Hexaethylene glycol monododecyl ...	450	C24H50O7	003055-96-7	38
5	1H-Pyrimidine-2,4-dione, 1-(2-et...	232	C9H13FN2O4	1000310-13-7	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071017\
Data File : BF096597.D
Acq On : 10 Jul 2017 13:29
Operator : SJ/JU
Sample : I4086-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

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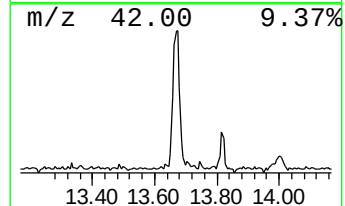
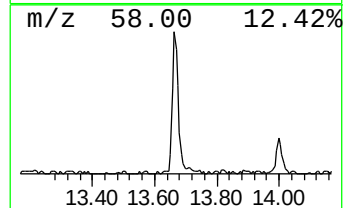
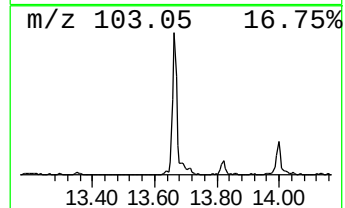
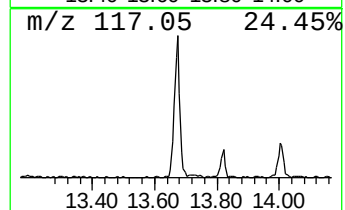
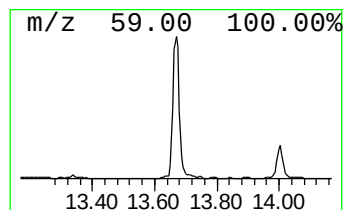
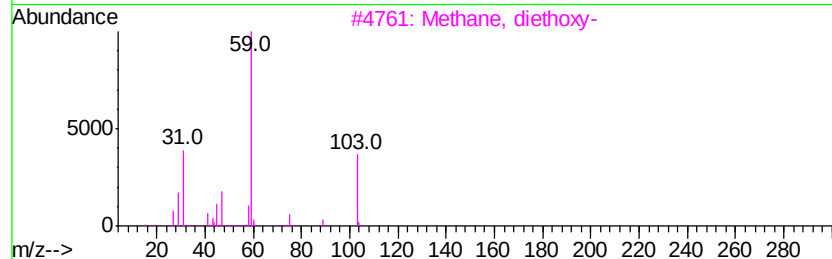
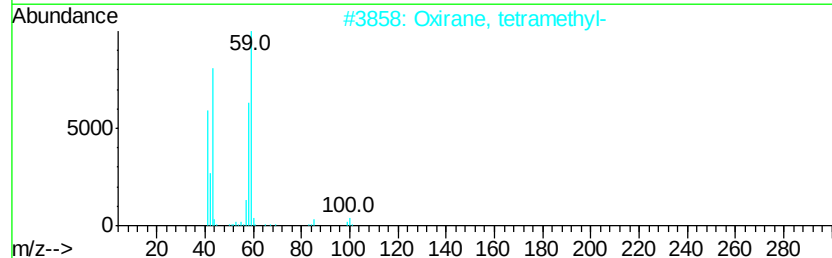
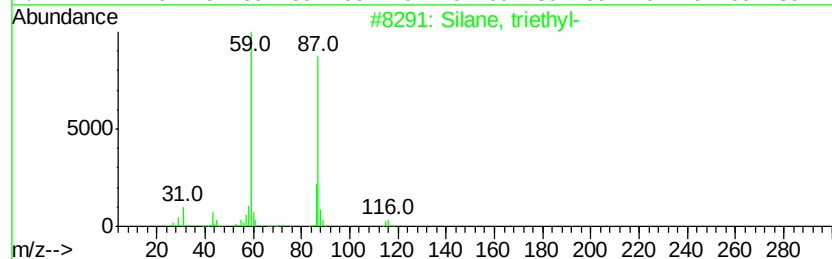
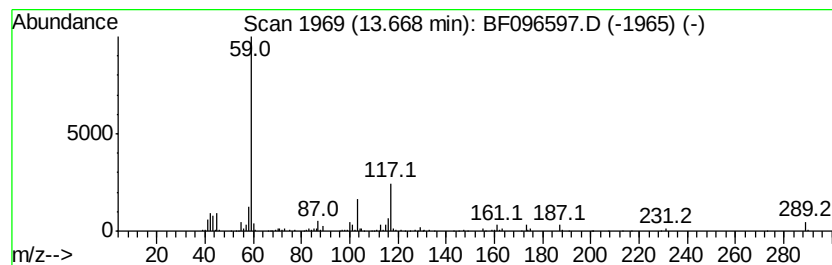
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown13.67 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	13.79 ng	740869	Chrysene-d12	13.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Silane, triethyl-	116	C6H16Si	000617-86-7	47
2		Oxirane, tetramethyl-	100	C6H12O	005076-20-0	47
3		Methane, diethoxy-	104	C5H12O2	000462-95-3	47
4		Methyl 15-methoxyhexadecanoate	300	C18H36O3	1000110-17-9	47
5		Methyl 16-methoxyheptadecanoate	314	C19H38O3	1000110-18-2	47



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071017\
Data File : BF096597.D
Acq On : 10 Jul 2017 13:29
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Sample : I4086-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

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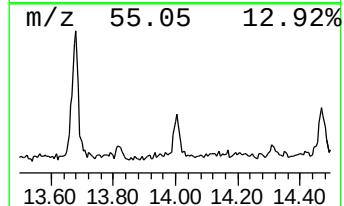
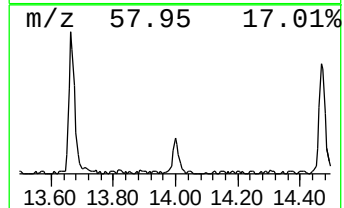
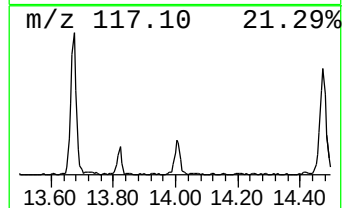
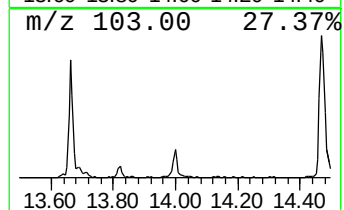
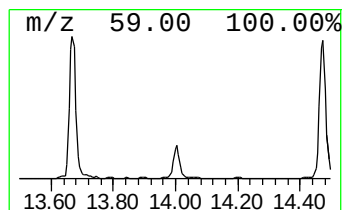
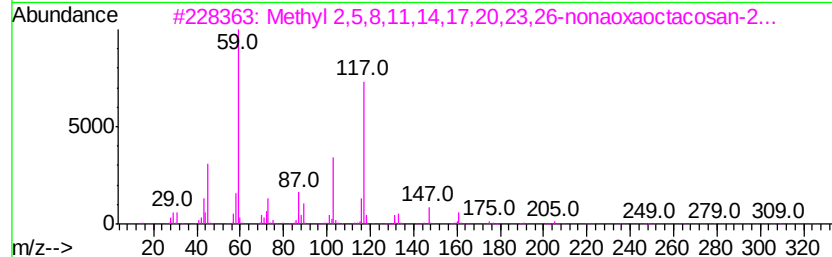
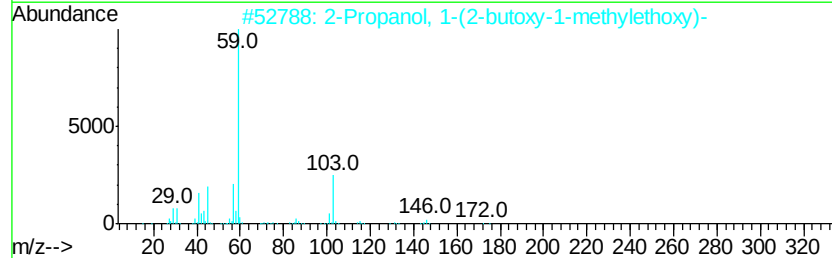
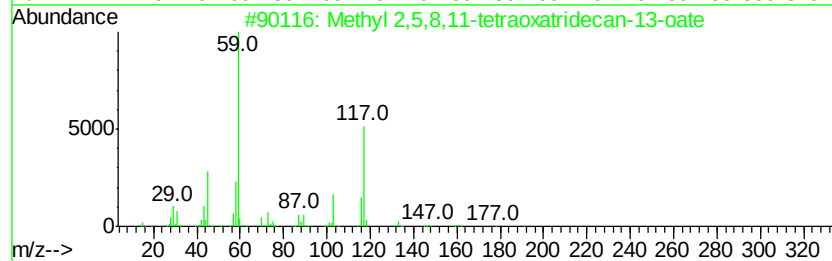
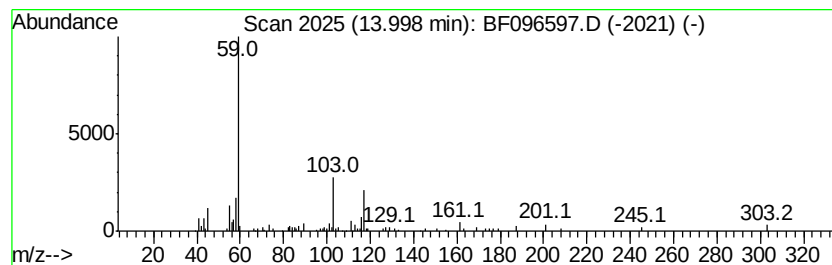
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Methyl 2,5,8,11-tetraoxatri... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.00	3.01 ng	161721	Chrysene-d12	13.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl 2,5,8,11-tetraoxatridecan...	236	C10H20O6	1000366-78-2	64
2		2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	50
3		Methyl 2,5,8,11,14,17,20,23,26-n...	456	C20H40O11	1000366-81-1	50
4		Methyl 2,5,8,11,14,17,20,23-octa...	412	C18H36O10	1000366-81-2	50
5		Butanamide, 3,3-dimethyl-	115	C6H13NO	000926-04-5	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071017\
Data File : BF096597.D
Acq On : 10 Jul 2017 13:29
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Sample : I4086-01
Misc :
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Instrument :
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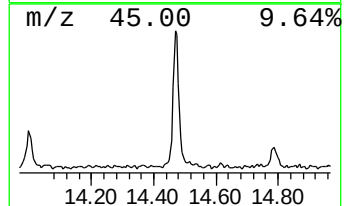
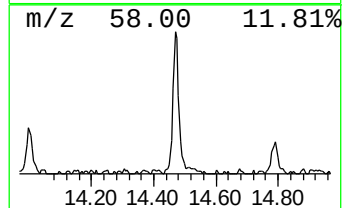
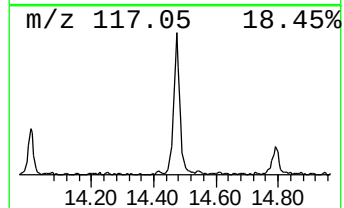
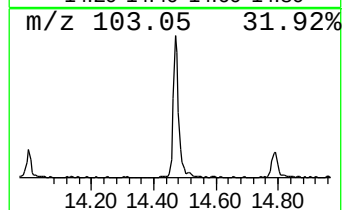
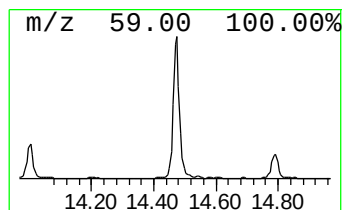
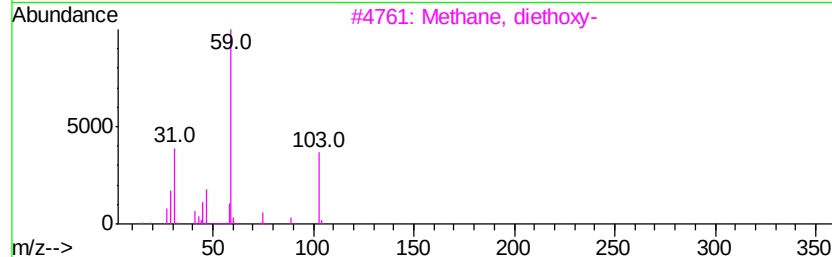
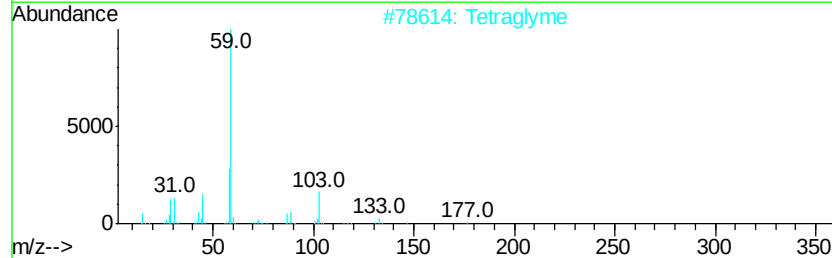
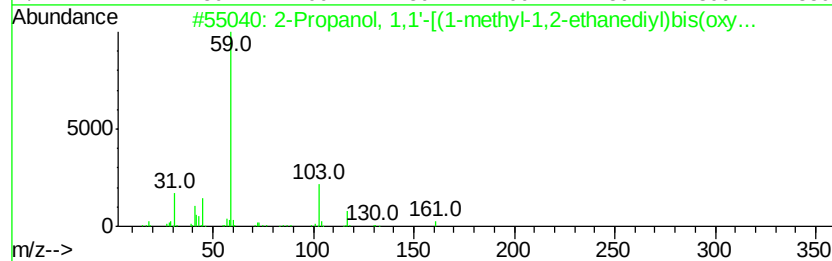
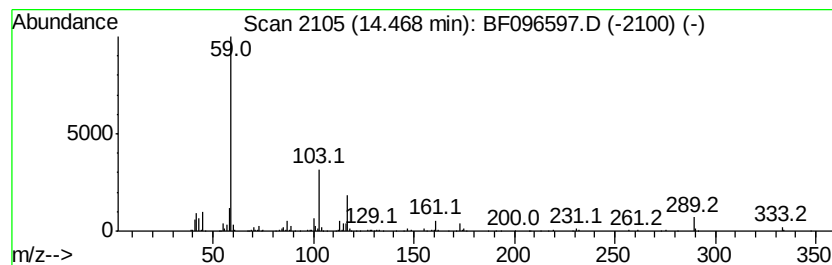
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 2-Propanol, 1,1'-[(1-methyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.47	11.28 ng	605966	Chrysene-d12	13.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	59
2		Tetraolyme	222	C10H22O5	000143-24-8	50
3		Methane, diethoxy-	104	C5H12O2	000462-95-3	50
4		Methyl 2,5,8,11,14,17,20,23,26,2...	500	C22H44O12	1000366-81-0	45
5		Propanol, [(butoxymethylethoxy)m...	248	C13H28O4	055934-93-5	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071017\
Data File : BF096597.D
Acq On : 10 Jul 2017 13:29
Operator : SJ/JU
Sample : I4086-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

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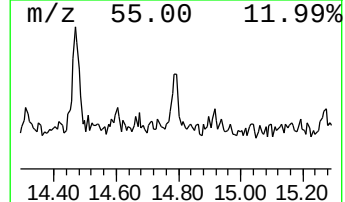
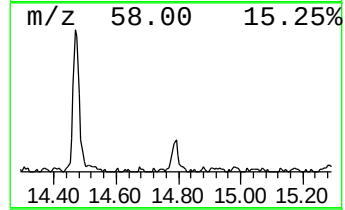
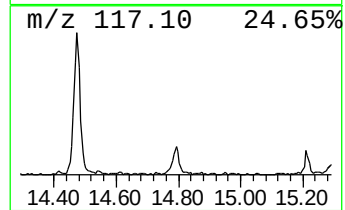
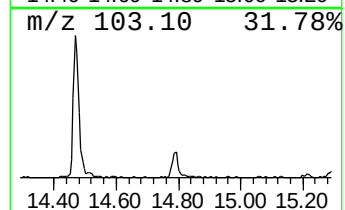
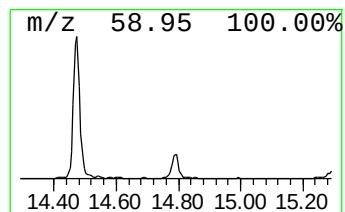
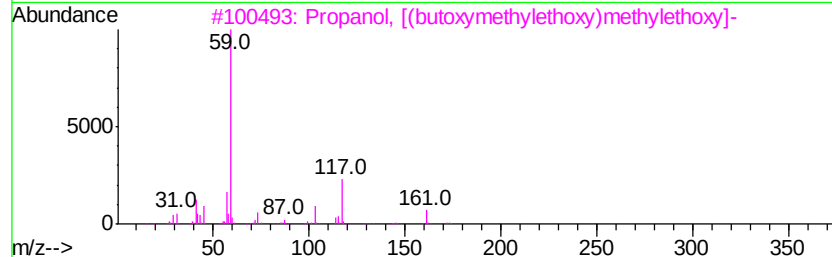
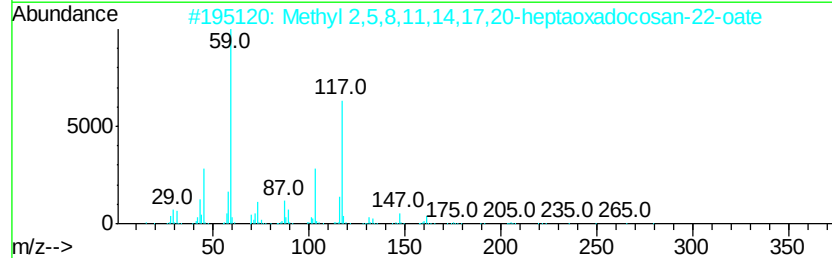
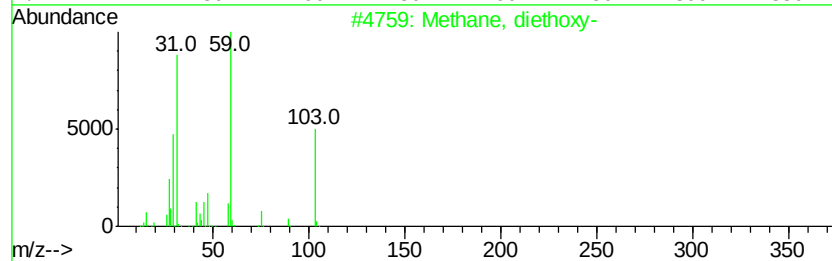
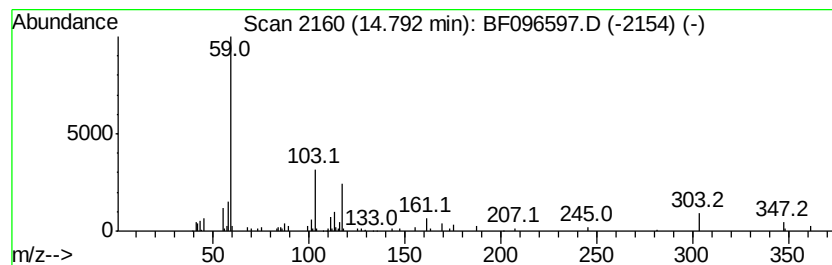
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown14.79 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.79	3.19 ng	121101	Perylene-d12	15.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methane, diethoxy-	104	C5H12O2	000462-95-3	47
2		Methyl 2,5,8,11,14,17,20-heptaox...	368	C16H32O9	1000366-81-3	45
3		Propanol, [(butoxymethyl)ethoxy]m...	248	C13H28O4	055934-93-5	42
4		Tri(propylene glycol) propyl ether	234	C12H26O4	096077-04-2	42
5		Hydrazine, 1-methyl-1-(2-methylp...	102	C5H14N2	020240-63-5	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071017\
Data File : BF096597.D
Acq On : 10 Jul 2017 13:29
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Sample : I4086-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

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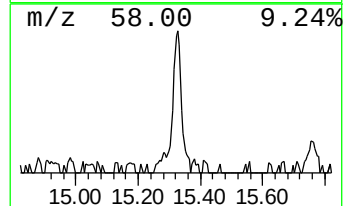
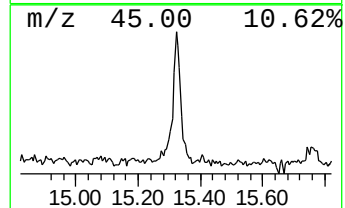
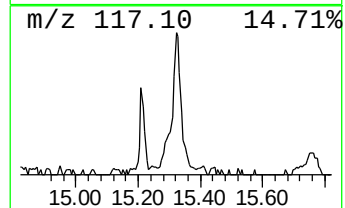
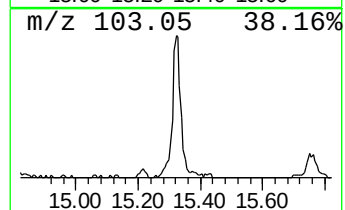
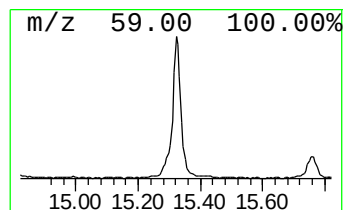
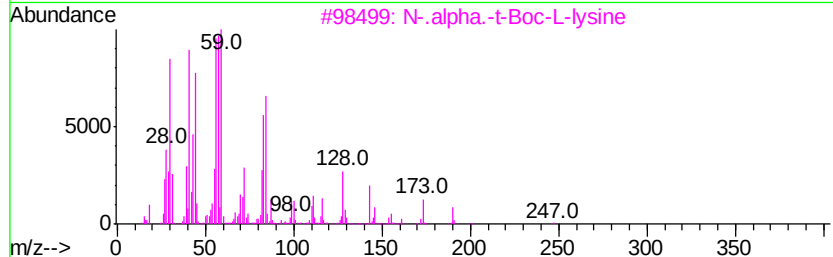
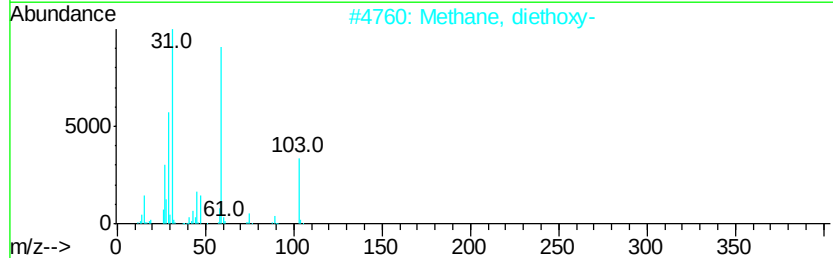
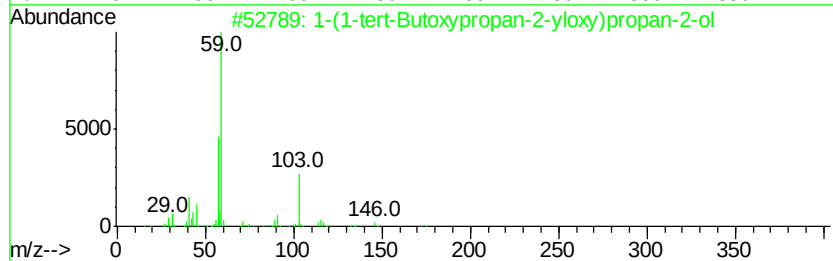
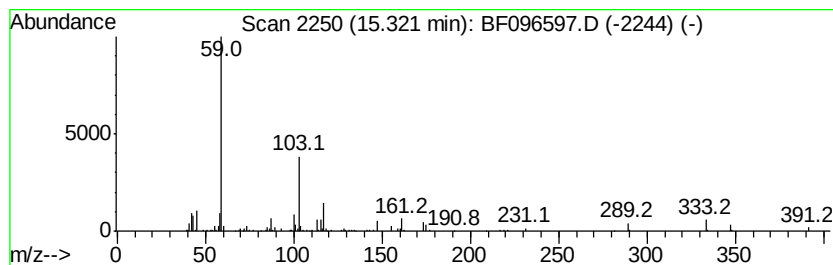
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-(1-tert-Butoxypropan-2-yl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.32	8.46 ng	321675	Perylene-d12	15.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-(1-tert-Butoxypropan-2-yloxy)p...	190	C10H22O3	132739-31-2	59
2		Methane, diethoxy-	104	C5H12O2	000462-95-3	53
3		N-.alpha.-t-Boc-L-lysine	246	C11H22N2O4	013734-28-6	43
4		4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	38
5		Methyl 2,5,8,11-tetraoxatridecan...	236	C10H20O6	1000366-78-2	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071017\
Data File : BF096597.D
Acq On : 10 Jul 2017 13:29
Operator : SJ/JU
Sample : I4086-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
A2882

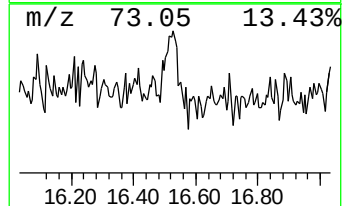
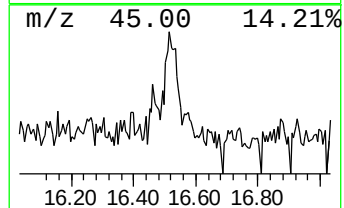
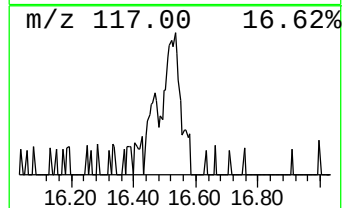
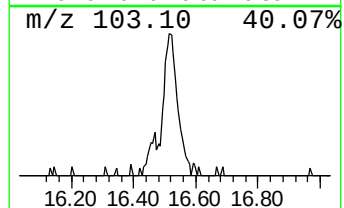
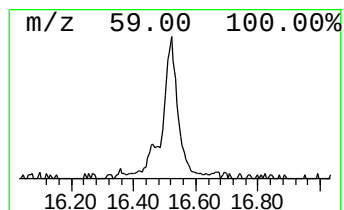
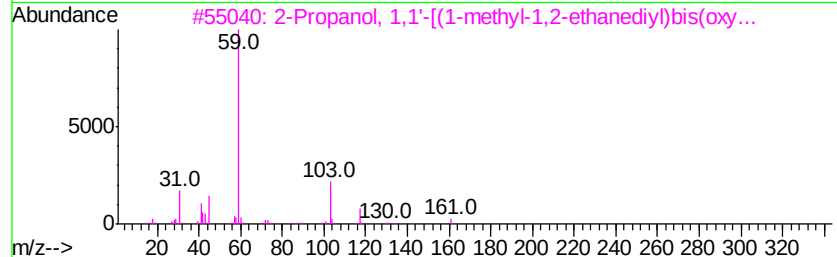
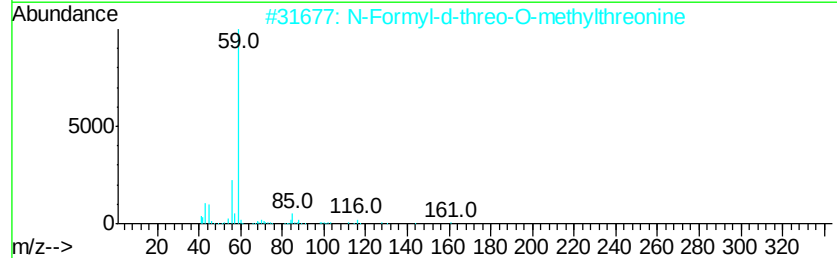
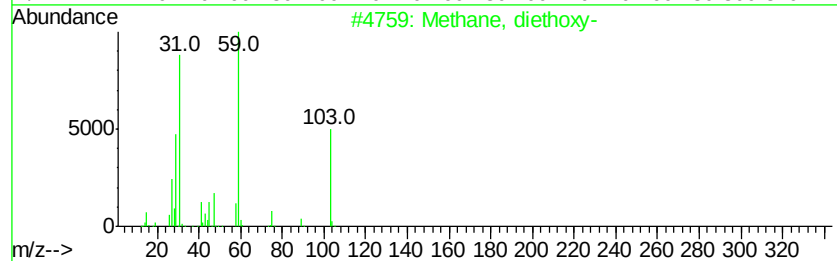
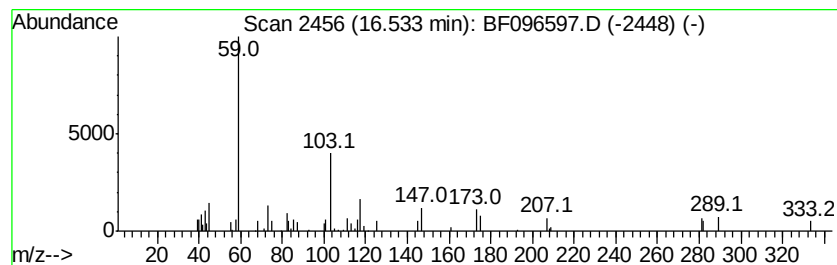
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown16.53 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.53	2.82 ng	107241	Perylene-d12	15.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methane, diethoxy-	104	C5H12O2	000462-95-3	40
2		N-Formyl-d-threo-O-methylthreonine	161	C6H11NO4	1000214-69-5	35
3		2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	33
4		Butanedioic acid, diazo-, dimeth...	172	C6H8N2O4	055514-36-8	33
5		(R)-3-Hydroxybutyric acid, methy...	132	C6H12O3	1000332-95-6	32



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071017\
Data File : BF096597.D
Acq On : 10 Jul 2017 13:29
Operator : SJ/JU
Sample : I4086-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
A2882

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.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
2-Pentanone, 4-hy...	4.86	6.5	ng	250834	1	6.68	766431	20.0
unknown6.45	6.45	72.9	ng	2795320	1	6.68	766431	20.0
Heptyl isobutyl c...	6.75	2.7	ng	102691	1	6.68	766431	20.0
unknown8.22	8.22	2.1	ng	113511	2	7.96	1080560	20.0
unknown11.73	11.73	2.9	ng	165824	4	11.19	1128330	20.0
unknown12.47	12.47	2.5	ng	142806	4	11.19	1128330	20.0
unknown13.67	13.67	13.8	ng	740869	5	13.82	1074620	20.0
Methyl 2,5,8,11-t...	14.00	3.0	ng	161721	5	13.82	1074620	20.0
2-Propanol, 1,1'-...	14.47	11.3	ng	605966	5	13.82	1074620	20.0
unknown14.79	14.79	3.2	ng	121101	6	15.22	760067	20.0
1-(1-tert-Butoxyp...	15.32	8.5	ng	321675	6	15.22	760067	20.0
unknown16.53	16.53	2.8	ng	107241	6	15.22	760067	20.0