

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122116\
 Data File : BF091863.D
 Acq On : 22 Dec 2016 4:50
 Operator : UM/SJ
 Sample : H6156-09
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-18

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-BF122116.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	2.213	9	11	39	rVB4	75100	442385	2.86%	0.464%
2	3.618	130	134	138	rBV	271654	455903	2.94%	0.478%
3	4.910	245	247	251	rBV	687695	873511	5.64%	0.916%
4	5.321	278	283	284	rBV	837417	1741246	11.24%	1.826%
5	5.344	284	285	288	rVB	3282740	2900015	18.72%	3.041%
6	5.573	302	305	307	rBV	2435883	3336140	21.53%	3.498%
7	5.904	331	334	336	rBV	278614	325235	2.10%	0.341%
8	6.384	374	376	378	rVV	1865155	2025390	13.07%	2.124%
9	6.430	378	380	381	rVV	588035	590707	3.81%	0.619%
10	6.453	381	382	385	rVB	303013	252483	1.63%	0.265%
11	6.510	385	387	389	rBV	4249945	5591563	36.09%	5.863%
12	6.567	389	392	396	rVB2	357528	699731	4.52%	0.734%
13	6.739	405	407	410	rBV	1171621	1600957	10.33%	1.679%
14	6.807	410	413	415	rVV	729133	827634	5.34%	0.868%
15	6.899	419	421	426	rVB2	4281628	5729566	36.98%	6.007%
16	7.013	428	431	433	rVV	5891576	7380808	47.64%	7.739%
17	7.310	451	457	460	rBV	3475199	4482041	28.93%	4.699%
18	7.779	495	498	500	rVV2	1248462	1947029	12.57%	2.041%
19	7.813	500	501	505	rVV	1424738	2030794	13.11%	2.129%
20	7.905	507	509	511	rVB	170285	169532	1.09%	0.178%
21	8.065	517	523	524	rBV2	8679466	15494121	100.00%	16.245%
22	8.099	524	526	529	rVB	842105	1080455	6.97%	1.133%
23	8.533	563	564	570	rVV2	156885	265901	1.72%	0.279%
24	8.693	573	578	581	rVV3	147288	215179	1.39%	0.226%
25	8.842	588	591	593	rBV	2416916	2338355	15.09%	2.452%
26	8.888	593	595	598	rVV	210867	227257	1.47%	0.238%
27	9.116	611	615	617	rBV	5432566	7991587	51.58%	8.379%
28	9.448	641	644	646	rBV	120815	169154	1.09%	0.177%
29	9.642	658	661	663	rBV	1500094	1309906	8.45%	1.373%
30	9.779	670	673	675	rBV	2528245	2738467	17.67%	2.871%
31	9.813	675	676	678	rVB	708899	534694	3.45%	0.561%
32	10.031	693	695	698	rBV	385635	422774	2.73%	0.443%
33	10.568	739	742	745	rBV2	3509966	4531773	29.25%	4.751%
34	11.265	800	803	805	rBV	2615303	2650172	17.10%	2.779%

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35	12.854	939	942	944	rBV	6210745	6928085	44.71%	7.264%
36	13.894	1030	1033	1036	rBV	2368080	2268086	14.64%	2.378%
37	15.288	1152	1155	1157	rBV	1449185	1800753	11.62%	1.888%
38	15.334	1157	1159	1168	rVB	119352	209966	1.36%	0.220%
39	15.723	1190	1193	1198	rBV	135458	209645	1.35%	0.220%
40	16.180	1230	1233	1239	rVB	137981	248177	1.60%	0.260%
41	16.717	1276	1280	1284	rBV	84740	177299	1.14%	0.186%
42	17.346	1331	1335	1342	rBV	63499	162670	1.05%	0.171%

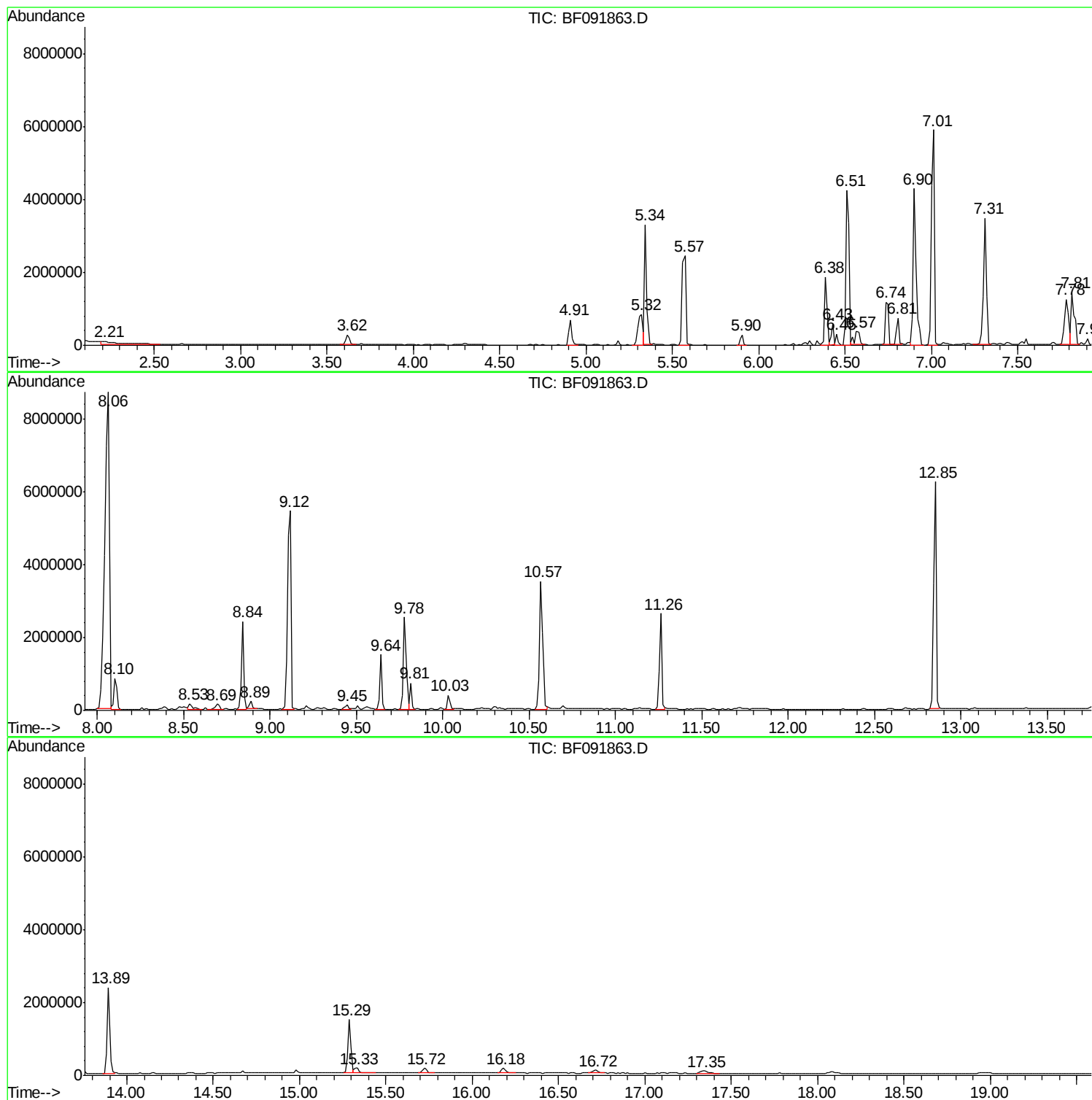
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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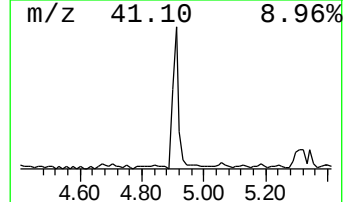
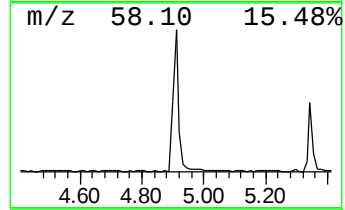
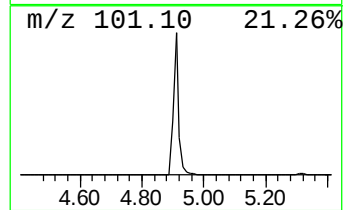
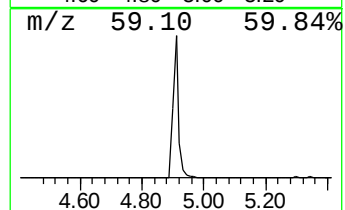
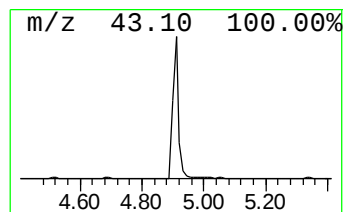
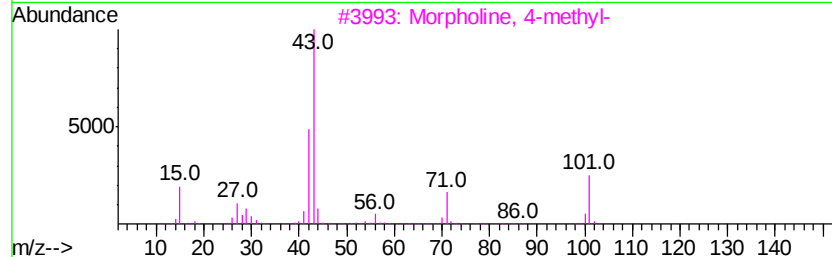
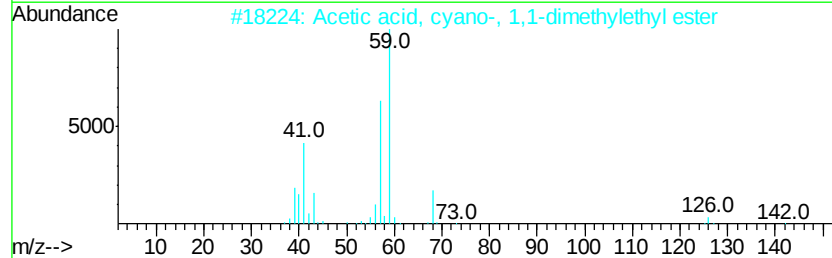
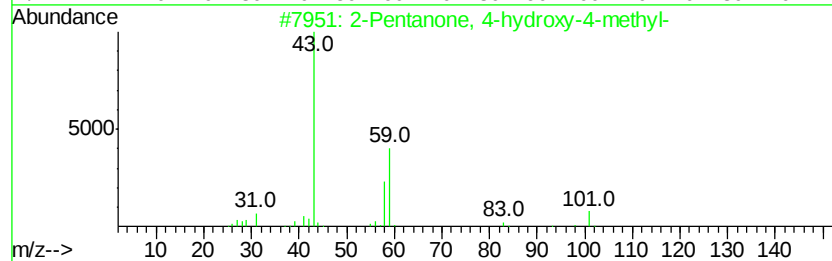
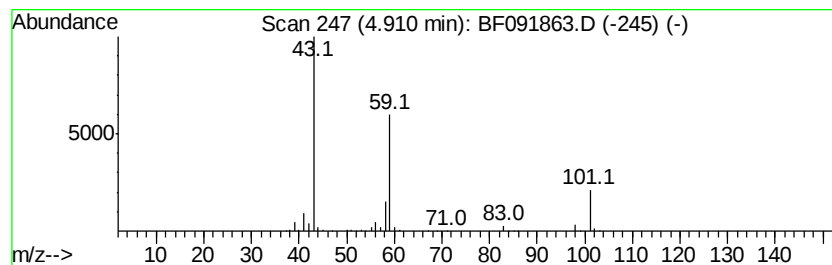
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.91	10.91 ng	873511	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	9



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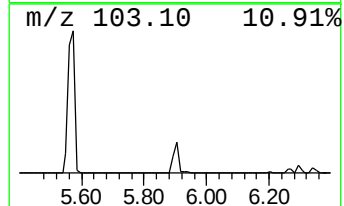
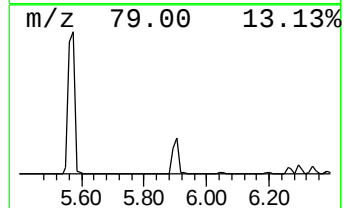
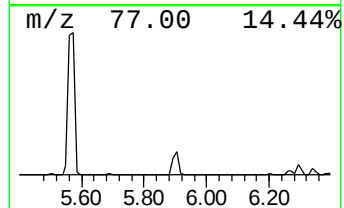
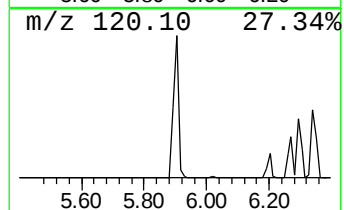
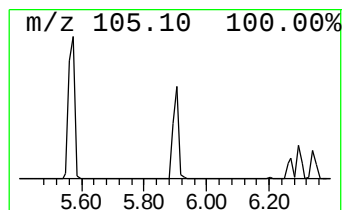
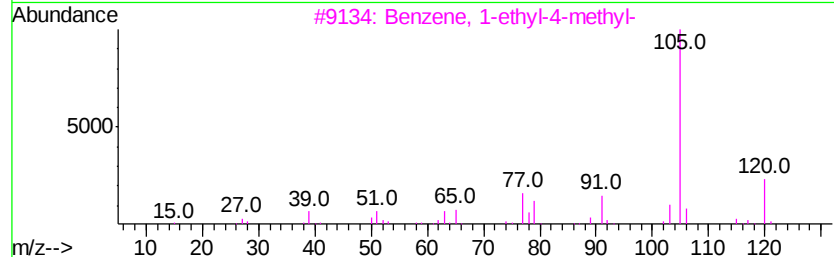
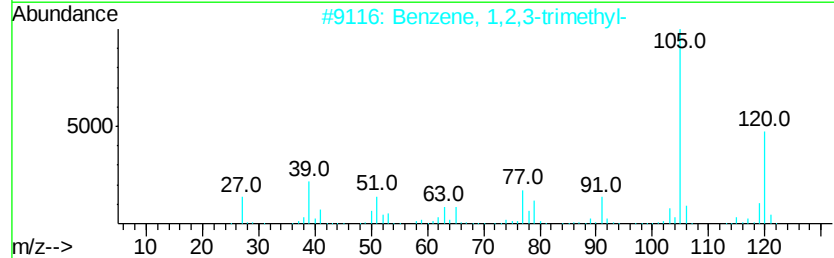
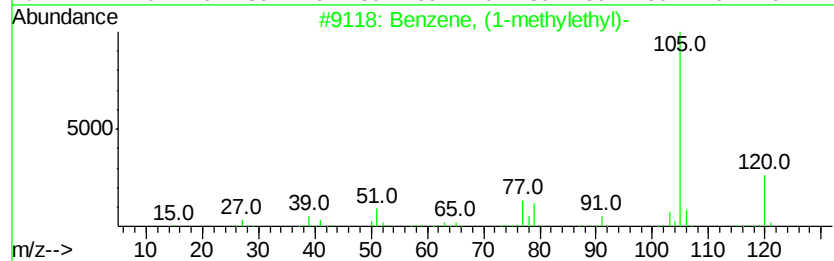
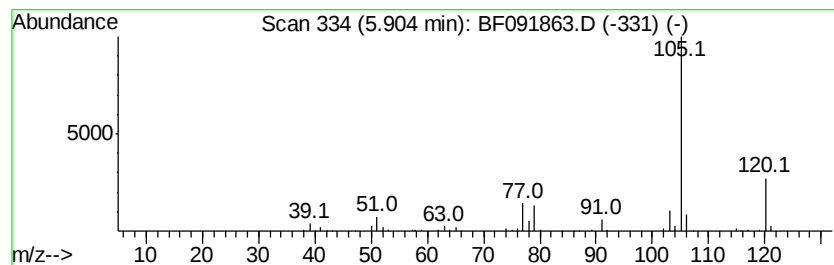
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Benzene, (1-methylethyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.90	4.06 ng	325235	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	80
4		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	78
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	72



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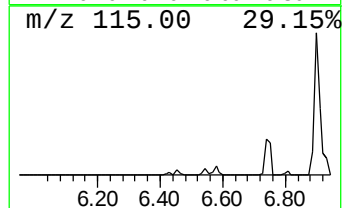
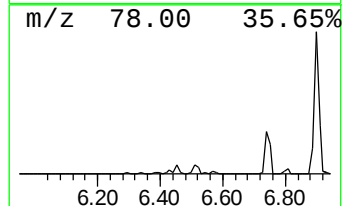
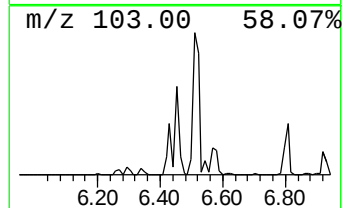
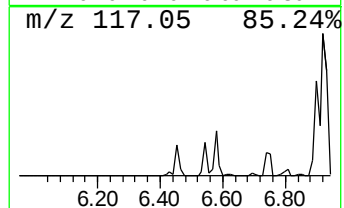
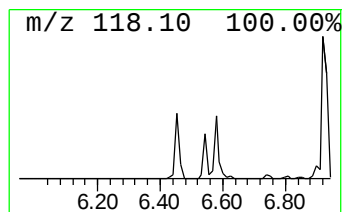
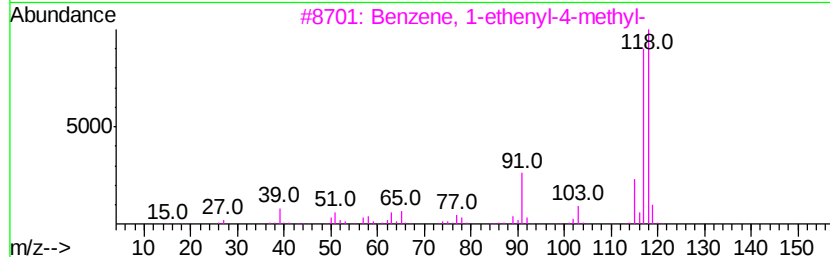
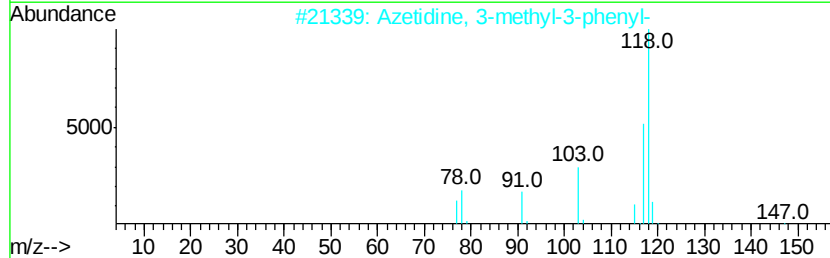
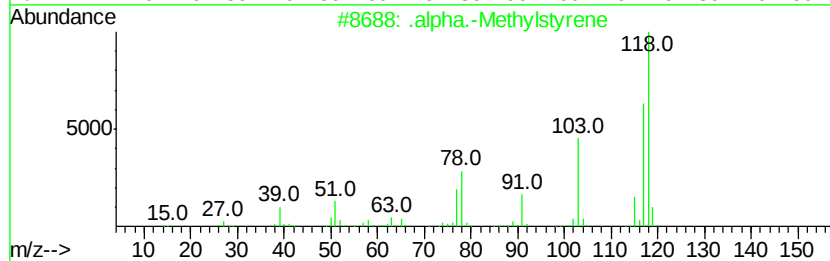
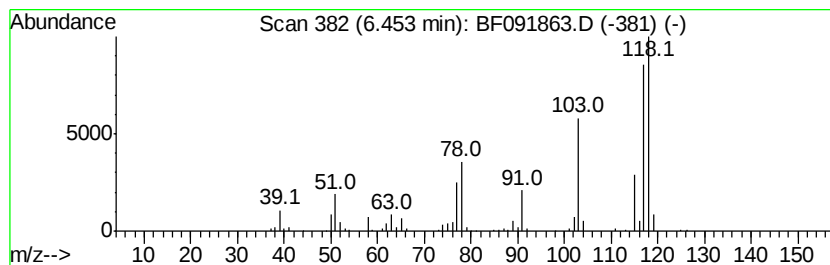
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 .alpha.-Methylstyrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.45	3.15 ng	252483	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.alpha.-Methylstvrene	118	C9H10	000098-83-9	94
2		Azetidine, 3-methyl-3-phenyl-	147	C10H13N	005961-33-1	64
3		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	53
4		Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	49
5		Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	43



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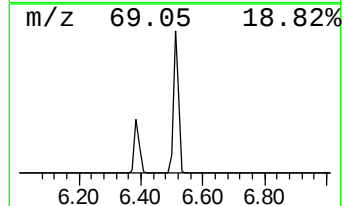
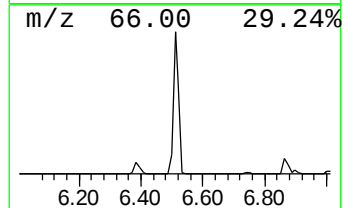
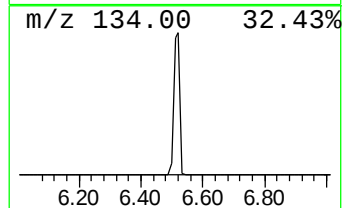
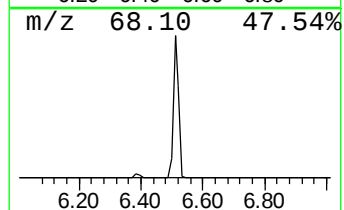
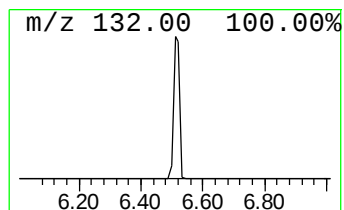
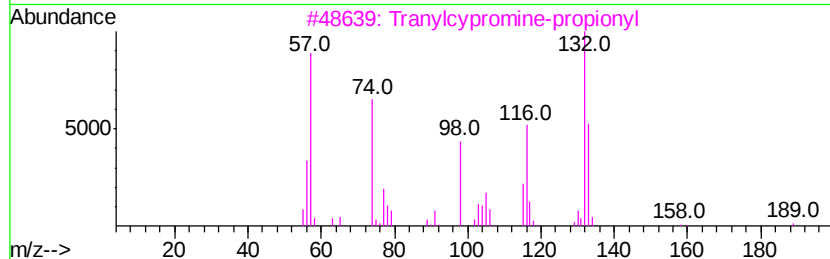
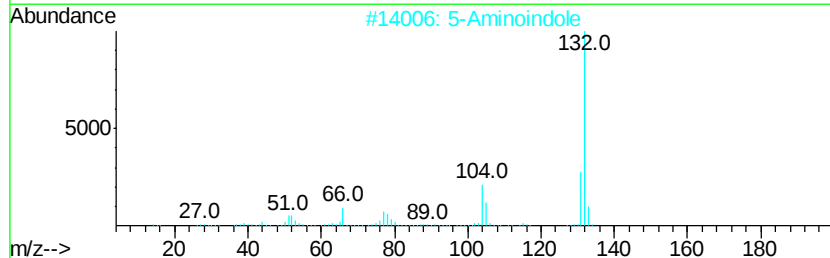
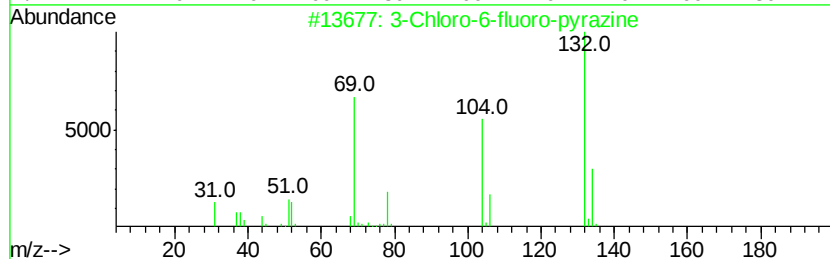
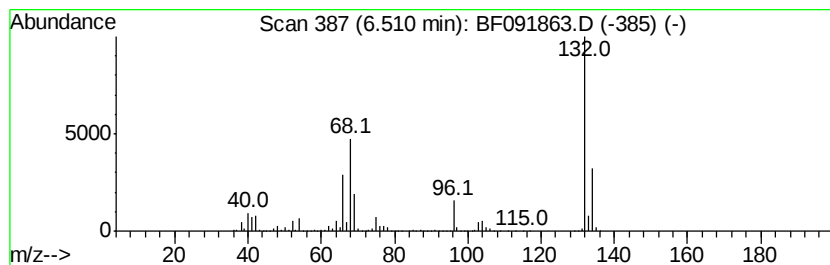
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Peak Number 7 unknown6.51 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	69.85 ng	5591560	1,4-Dichlorobenzene-d4	6.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5		Aminoindole	132	C8H8N2	005192-03-0	12
3	Tranvlcv		promine-propionyl	189	C12H15NO	1000123-86-3	10
4	Benzene,		1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5	Bicyclo[4.2.0]octa-1,3,5-triene,		...	132	C10H12	028749-81-7	9



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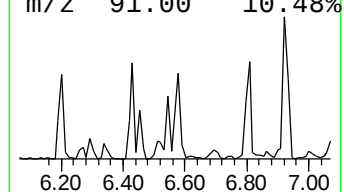
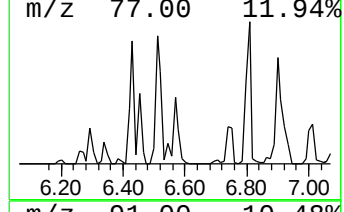
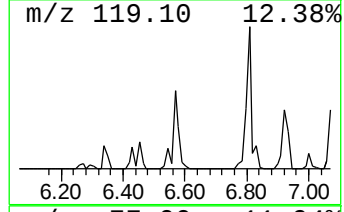
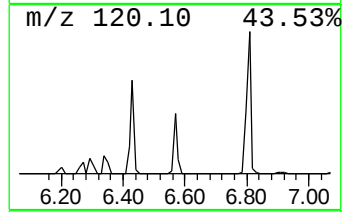
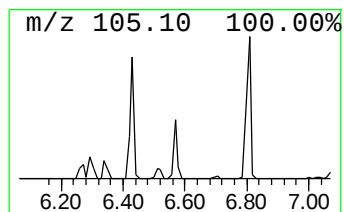
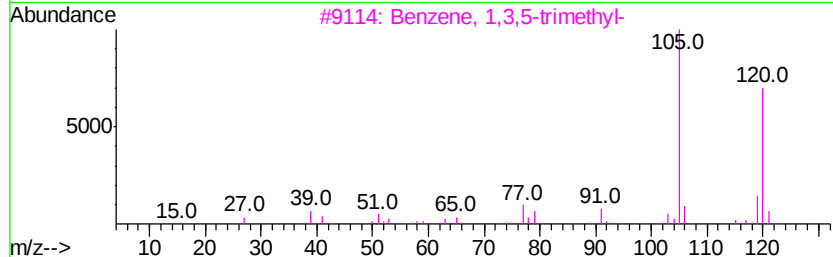
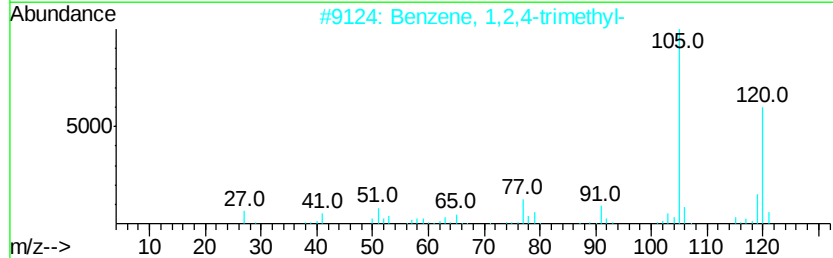
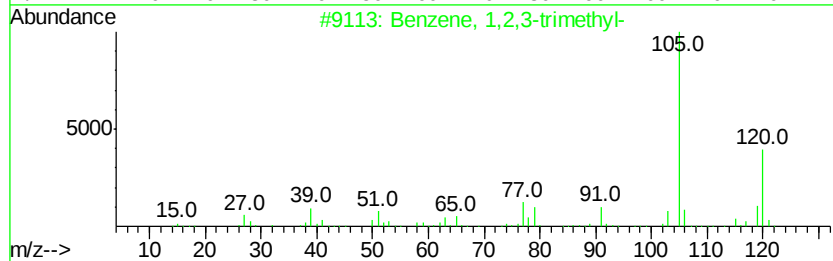
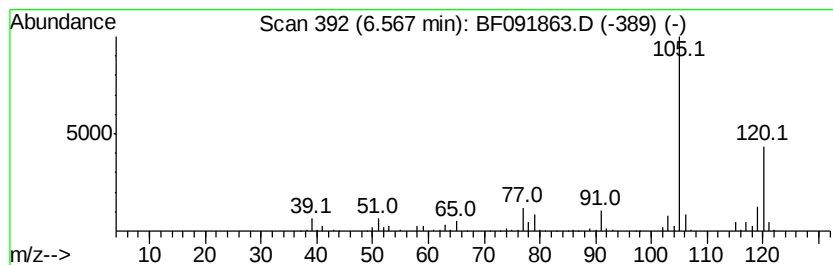
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	8.74 ng	699731	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122116\
Data File : BF091863.D
Acq On : 22 Dec 2016 4:50
Operator : UM/SJ
Sample : H6156-09
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-18

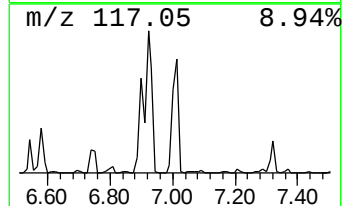
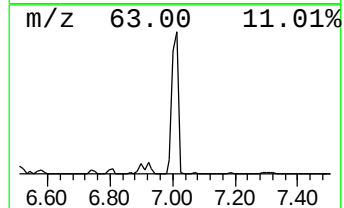
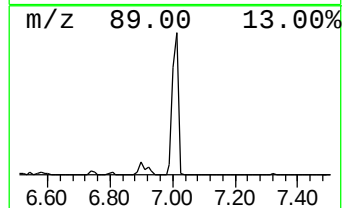
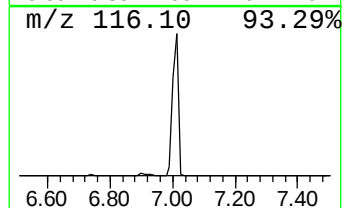
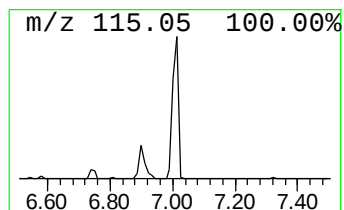
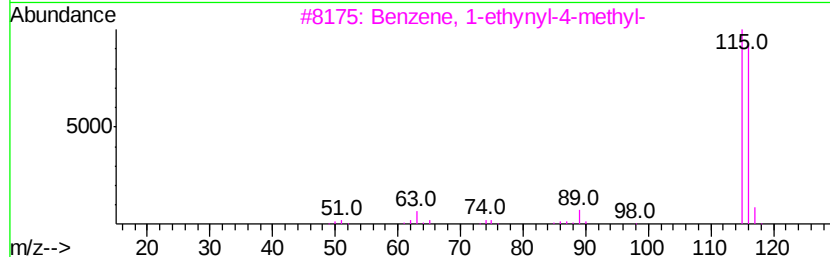
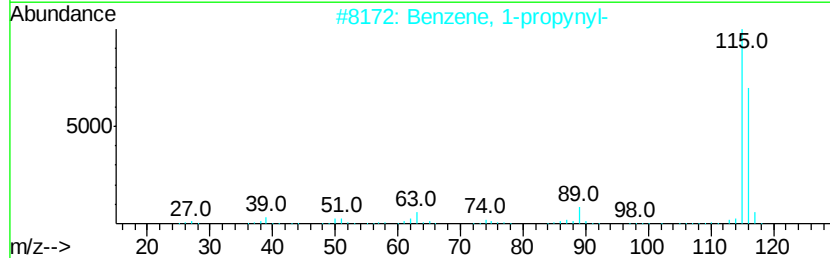
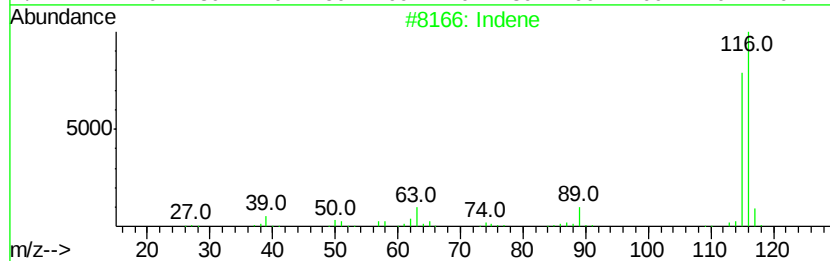
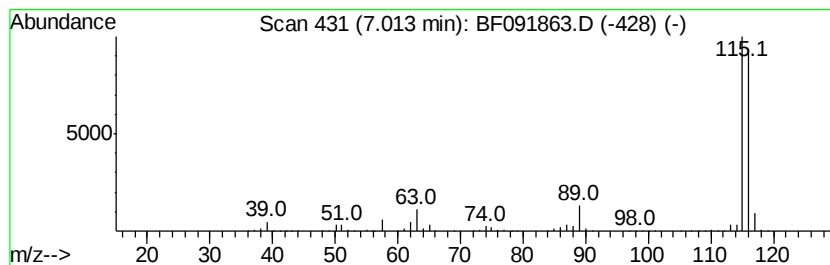
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Indene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.01	92.20 ng	7380810	1,4-Dichlorobenzene-d4	6.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	97
2			Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
3			Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91
4			Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91
5			Benzene,1-ethynyl-2-methyl-	116	C9H8	000766-47-2	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122116\
Data File : BF091863.D
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Sample : H6156-09
Misc :
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Instrument :
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ClientSampleId :
MW-18

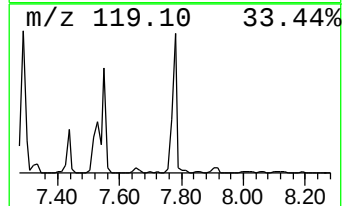
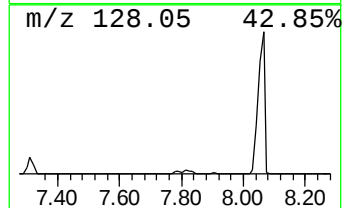
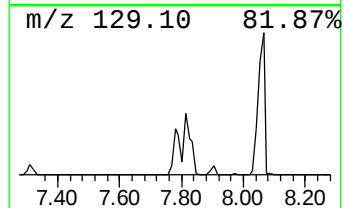
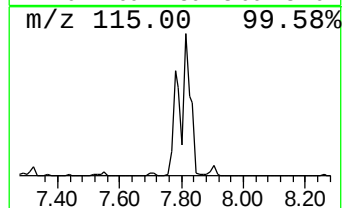
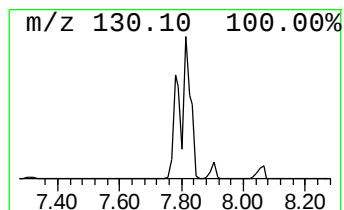
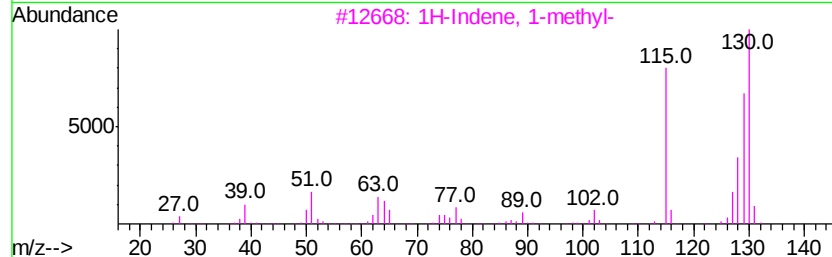
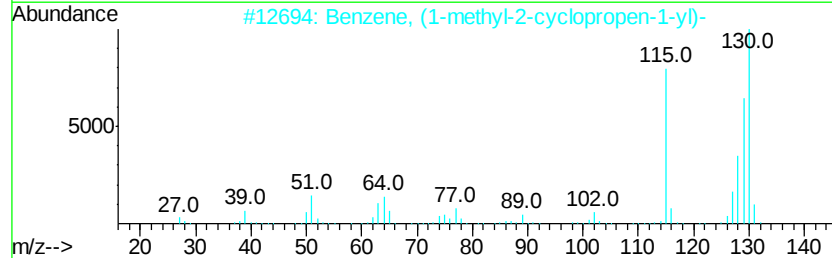
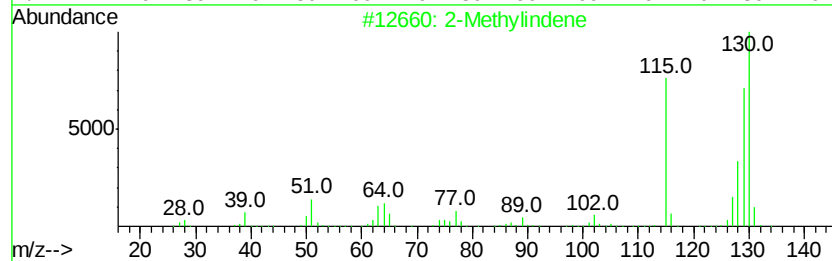
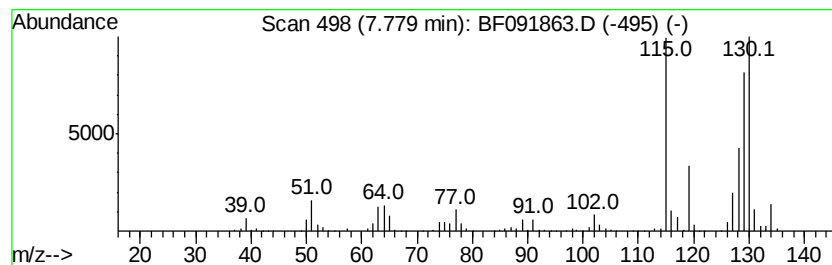
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2-Methylindene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.78	2.51 ng	1947030	Naphthalene-d8	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Methylindene	130	C10H10	002177-47-1	97
2		Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	97
3		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
4		Benzene, 1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	96
5		1H-Indene, 3-methyl-	130	C10H10	000767-60-2	94



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
Data File : BF091863.D
Acq On : 22 Dec 2016 4:50
Operator : UM/SJ
Sample : H6156-09
Misc :
ALS Vial : 33 Sample Multiplier: 1

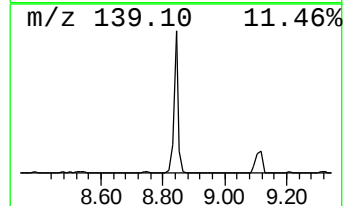
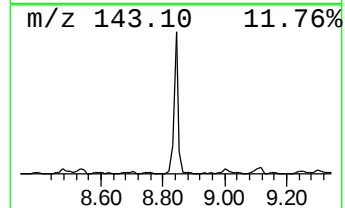
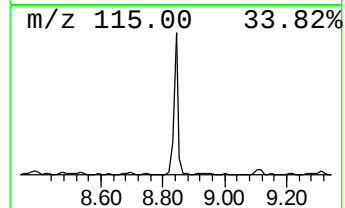
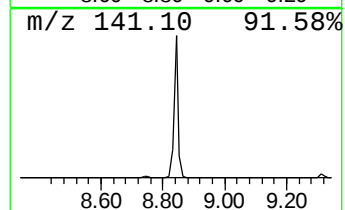
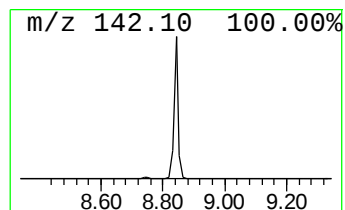
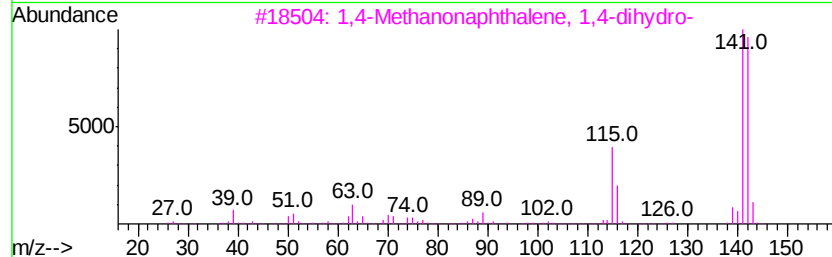
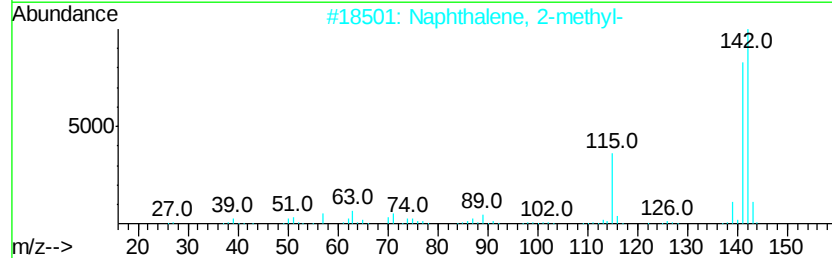
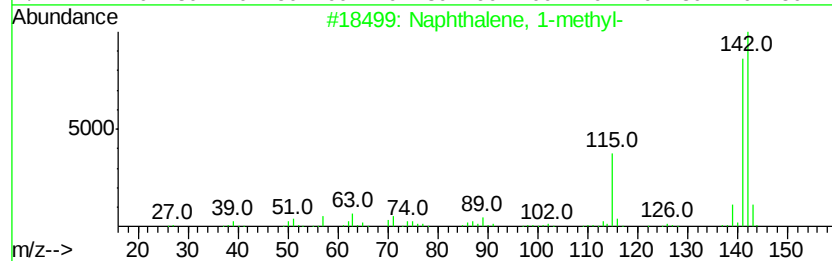
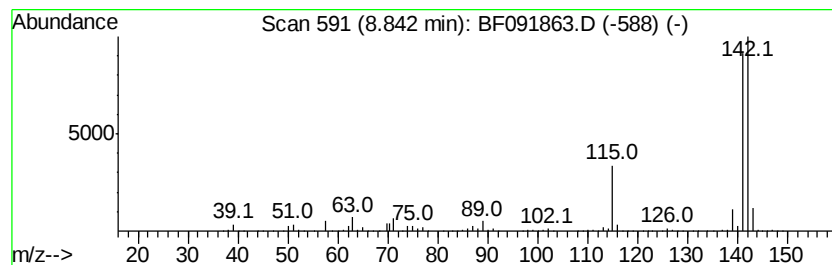
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Naphthalene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.84	3.02 ng	2338360	Naphthalene-d8	8.03
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 96
2		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
3		1,4-Methanonaphthalene, 1,4-dihv...	142 C11H10	004453-90-1 91
4		Benzocycloheptatriene	142 C11H10	000264-09-5 91
5		1H-Indene, 1-ethylidene-	142 C11H10	002471-83-2 64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
Data File : BF091863.D
Acq On : 22 Dec 2016 4:50
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Sample : H6156-09
Misc :
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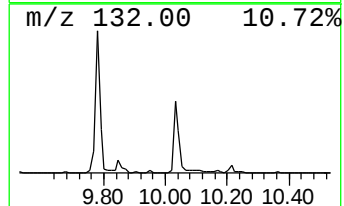
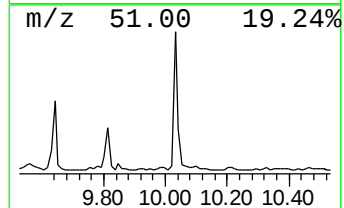
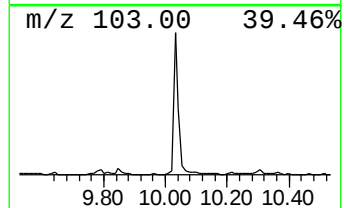
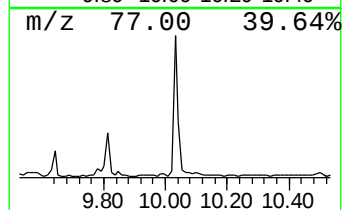
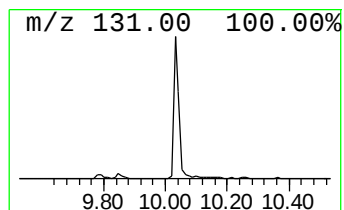
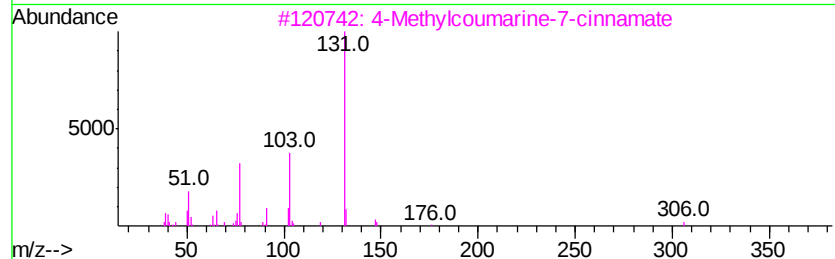
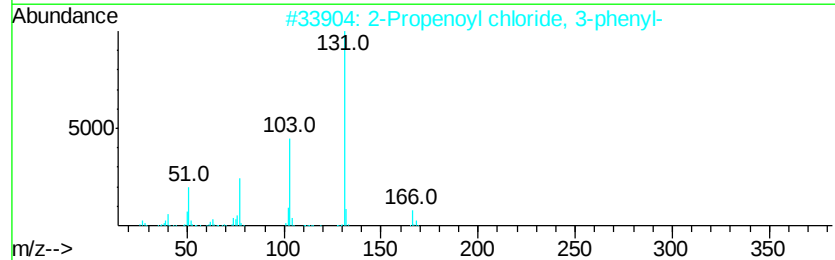
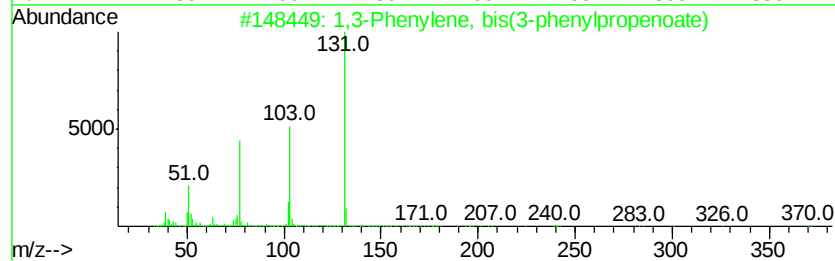
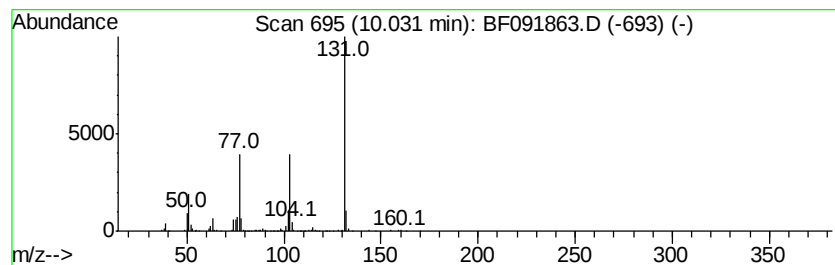
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 1,3-Phenylene, bis(3-phenyl... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.03	3.09 ng	422774	Acenaphthene-d10	9.78
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1,3-Phenylene, bis(3-phenylprope...	370 C24H18O4	1000259-62-4	90
2	2-Propenoyl chloride, 3-phenyl-	166 C9H7ClO	000102-92-1	86
3	4-Methylcoumarine-7-cinnamate	306 C19H14O4	300829-05-4	83
4	Propanedioic acid, nitrile, hydr...	229 C12H11N3O2	294878-35-6	78
5	Vinyl trans-cinnamate	174 C11H10O2	017719-70-9	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122116\
Data File : BF091863.D
Acq On : 22 Dec 2016 4:50
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Sample : H6156-09
Misc :
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Instrument :
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ClientSampleId :
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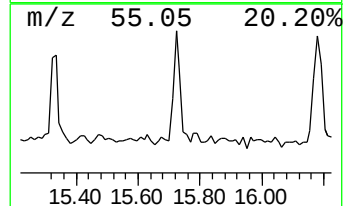
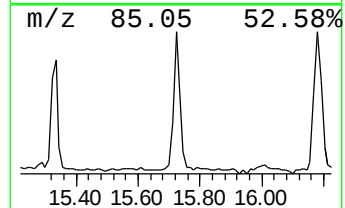
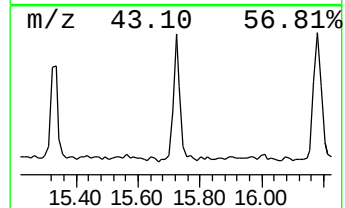
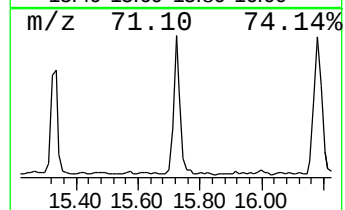
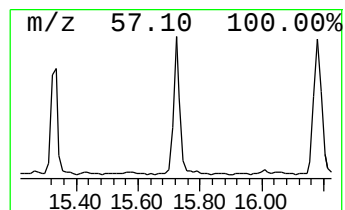
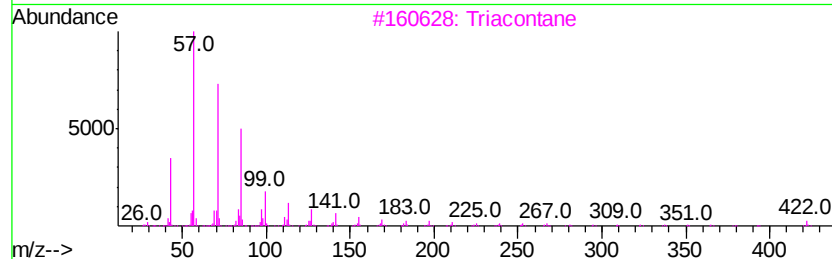
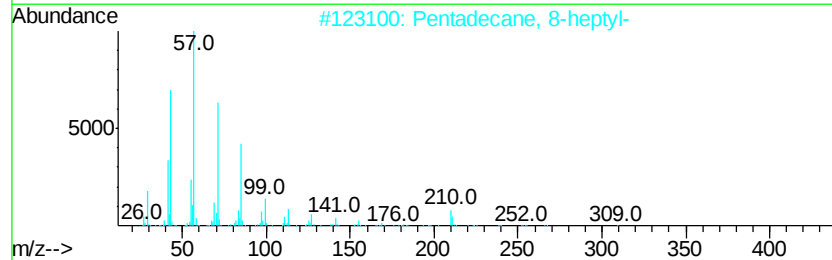
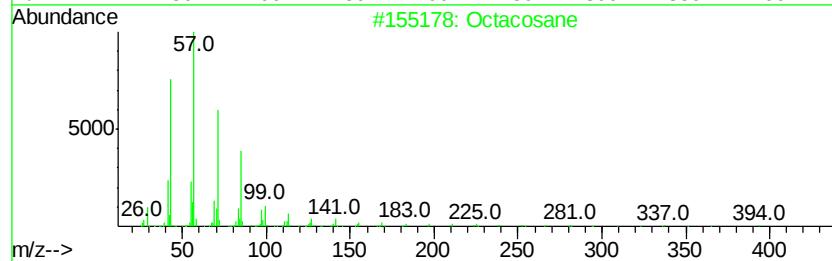
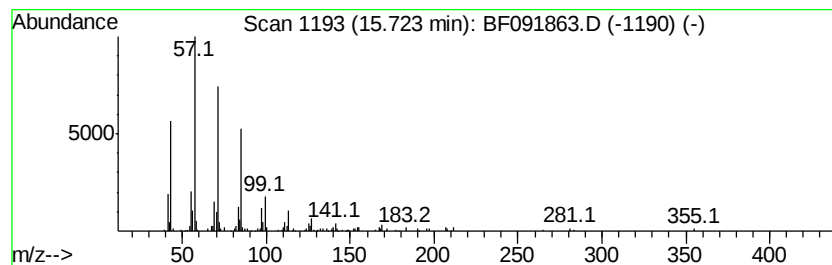
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Octacosane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.72	2.33 ng	209645	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
3		Triacontane	422	C30H62	000638-68-6	90
4		Nonacosane	408	C29H60	000630-03-5	90
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90



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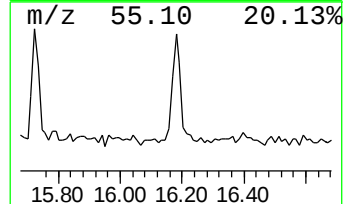
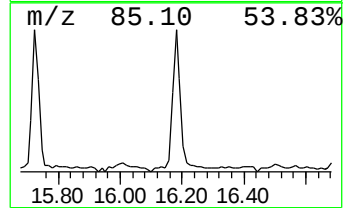
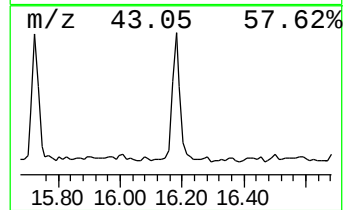
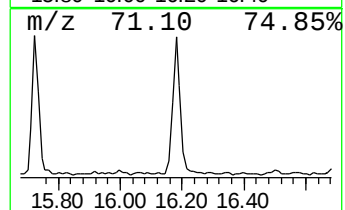
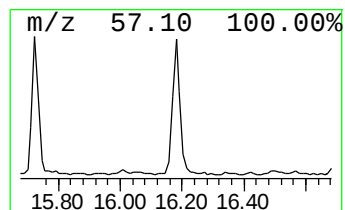
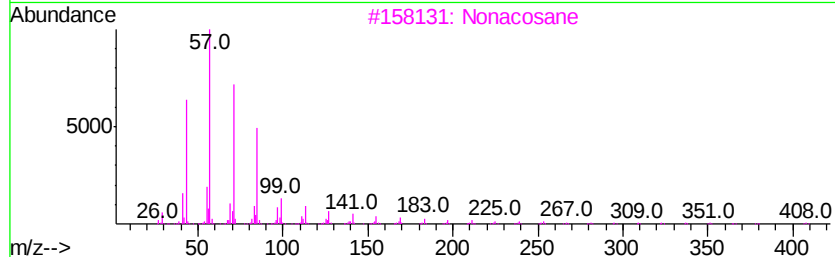
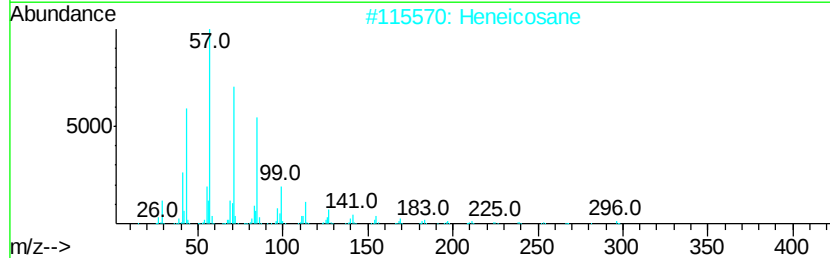
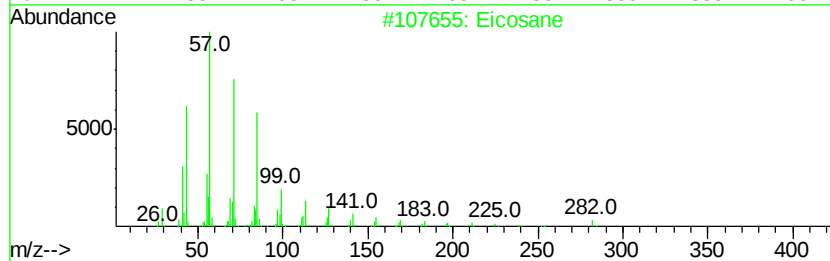
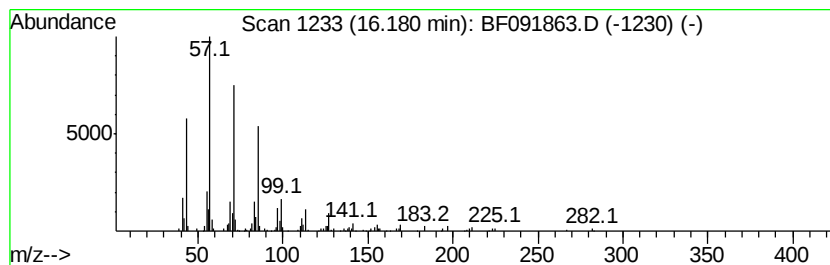
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Eicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.18	2.76 ng	248177	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Nonacosane	408	C29H60	000630-03-5	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122116\
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Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.91	10.9	ng	873511	1	6.75	1600960	20.0
Benzene, (1-methy...	5.90	4.1	ng	325235	1	6.75	1600960	20.0
.alpha.-Methylsty...	6.45	3.1	ng	252483	1	6.75	1600960	20.0
unknown6.51	6.51	69.8	ng	5591560	1	6.75	1600960	20.0
Benzene, 1,2,3-tr...	6.57	8.7	ng	699731	1	6.75	1600960	20.0
Indene	7.01	92.2	ng	7380810	1	6.75	1600960	20.0
2-Methylindene	7.78	2.5	ng	1947030	2	8.03	15494100	20.0
Naphthalene, 1-me...	8.84	3.0	ng	2338360	2	8.03	15494100	20.0
1,3-Phenylene, bi...	10.03	3.1	ng	422774	3	9.78	2738470	20.0
Octacosane	15.72	2.3	ng	209645	6	15.29	1800750	20.0
Eicosane	16.18	2.8	ng	248177	6	15.29	1800750	20.0