

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG052815\
 Data File : BG017324.D
 Acq On : 28 May 2015 14:30
 Operator : TP/IZ
 Sample : G2387-09MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-18-D13MS

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_G\METHODS\8270-BG051315.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	3.116	68	73	78	rBV	44890	58419	2.67%	0.099%
2	3.434	122	127	137	rBV	87419	126087	5.76%	0.214%
3	3.510	137	140	152	rVV	126553	188970	8.63%	0.321%
4	4.767	347	354	362	rBV	105502	156681	7.15%	0.266%
5	5.249	430	436	447	rBV	544924	793766	36.25%	1.350%
6	6.865	704	711	723	rVV2	598295	1237172	56.49%	2.103%
7	6.988	723	732	742	rVV	151943	255954	11.69%	0.435%
8	7.088	744	749	758	rVV	134273	197580	9.02%	0.336%
9	7.200	761	768	772	rVV	581578	978024	44.66%	1.663%
10	7.229	772	773	785	rVB	216623	221006	10.09%	0.376%
11	7.540	820	826	833	rVB	200524	303415	13.85%	0.516%
12	7.658	840	846	848	rBV	95872	151091	6.90%	0.257%
13	7.687	848	851	862	rVB	199967	308033	14.07%	0.524%
14	7.905	881	888	894	rBV2	240686	372535	17.01%	0.633%
15	7.975	894	900	903	rVV	377299	627495	28.65%	1.067%
16	8.005	903	905	913	rVB	211880	273324	12.48%	0.465%
17	8.128	918	926	931	rBV	259283	394777	18.03%	0.671%
18	8.199	931	938	944	rVB	144002	290540	13.27%	0.494%
19	8.463	974	983	993	rBV3	457097	1113404	50.84%	1.893%
20	8.727	1022	1028	1034	rVB	154918	242877	11.09%	0.413%
21	8.821	1036	1044	1048	rVV	335545	584898	26.71%	0.994%
22	8.862	1048	1051	1057	rVB	189598	264013	12.06%	0.449%
23	9.068	1079	1086	1097	rVB	65331	116456	5.32%	0.198%
24	9.380	1132	1139	1146	rVB	267285	423543	19.34%	0.720%
25	9.568	1163	1171	1178	rBV	170557	267195	12.20%	0.454%
26	9.650	1178	1185	1194	rVB	283912	453831	20.72%	0.772%
27	9.773	1197	1206	1211	rBV3	30493	67099	3.06%	0.114%
28	9.867	1217	1222	1230	rVB	176932	272579	12.45%	0.463%
29	10.114	1258	1264	1273	rBV	244670	392043	17.90%	0.667%
30	10.314	1289	1298	1305	rVB	243593	380929	17.39%	0.648%
31	10.449	1315	1321	1325	rBV	127488	211849	9.67%	0.360%
32	10.496	1325	1329	1336	rVB	288450	458772	20.95%	0.780%
33	10.619	1343	1350	1357	rBV2	143958	241557	11.03%	0.411%
34	10.784	1372	1378	1390	rVB2	219160	347621	15.87%	0.591%

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35	11.395	1474	1482	1485	rBV	76746	155282	7.09%	0.264%
36	11.430	1485	1488	1504	rVB3	51904	175686	8.02%	0.299%
37	11.777	1540	1547	1556	rBV	300850	468884	21.41%	0.797%
38	12.012	1581	1587	1595	rVB	172680	278443	12.71%	0.473%
39	12.117	1598	1605	1612	rBV	356942	560496	25.59%	0.953%
40	12.341	1633	1643	1651	rVB	355976	566326	25.86%	0.963%
41	12.470	1659	1665	1676	rVB2	391296	960635	43.87%	1.633%
42	12.746	1705	1712	1716	rVV	313621	458539	20.94%	0.780%
43	12.793	1716	1720	1723	rVV	255382	372767	17.02%	0.634%
44	12.834	1723	1727	1735	rVV	292573	451233	20.60%	0.767%
45	12.934	1735	1744	1751	rVB	810174	1169071	53.38%	1.988%
46	13.146	1773	1780	1783	rBV	426143	637357	29.10%	1.084%
47	13.181	1783	1786	1794	rVB	345092	523690	23.91%	0.890%
48	13.404	1814	1824	1831	rVB	270948	407641	18.61%	0.693%
49	13.692	1867	1873	1882	rBV	194631	280208	12.80%	0.476%
50	13.780	1882	1888	1893	rVV	497321	677229	30.92%	1.151%
51	13.839	1893	1898	1903	rVV	215799	324955	14.84%	0.552%
52	13.904	1903	1909	1916	rVB	309325	439582	20.07%	0.747%
53	14.004	1916	1926	1928	rBV	218203	312825	14.28%	0.532%
54	14.039	1928	1932	1938	rVB	402513	595561	27.20%	1.013%
55	14.233	1960	1965	1973	rVB	205433	317997	14.52%	0.541%
56	14.315	1973	1979	1984	rBV2	187669	291974	13.33%	0.496%
57	14.380	1984	1990	1995	rVB	485969	668075	30.51%	1.136%
58	14.444	1995	2001	2008	rBV	213388	322606	14.73%	0.549%
59	14.585	2018	2025	2029	rBV	524236	806643	36.83%	1.371%
60	14.691	2037	2043	2045	rBV2	352787	517822	23.65%	0.880%
61	14.714	2045	2047	2053	rVB	465455	554459	25.32%	0.943%
62	14.867	2068	2073	2082	rBV	350476	482675	22.04%	0.821%
63	14.949	2082	2087	2094	rVV	375777	501658	22.91%	0.853%
64	15.149	2112	2121	2126	rBV	525771	683788	31.22%	1.163%
65	15.367	2151	2158	2161	rBV2	953863	1381525	63.08%	2.349%
66	15.402	2161	2164	2169	rVV	313225	448747	20.49%	0.763%
67	15.455	2169	2173	2179	rVB	246272	319910	14.61%	0.544%
68	15.584	2186	2195	2200	rBV	584053	779131	35.58%	1.325%
69	15.655	2200	2207	2214	rVB	448342	604699	27.61%	1.028%
70	15.813	2227	2234	2242	rBV	884126	1205576	55.05%	2.050%
71	16.254	2304	2309	2316	rVB	445738	591809	27.02%	1.006%

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72	16.365	2322	2328	2336	rBV	418891	583844	26.66%	0.993%
73	16.536	2351	2357	2369	rBV	570714	717029	32.74%	1.219%
74	16.724	2383	2389	2401	rBV	664933	918197	41.93%	1.561%
75	17.065	2440	2447	2450	rBV	280837	398252	18.19%	0.677%
76	17.106	2450	2454	2460	rVV	631514	856155	39.09%	1.456%
77	17.194	2461	2469	2476	rVV	617780	853590	38.98%	1.451%
78	17.476	2510	2517	2527	rBV	648067	893180	40.79%	1.519%
79	17.969	2595	2601	2607	rBV	80303	109252	4.99%	0.186%
80	18.034	2607	2612	2618	rVB	683086	850650	38.84%	1.446%
81	19.127	2792	2798	2806	rBV	780906	1011988	46.21%	1.721%
82	19.321	2825	2831	2837	rBV	481126	605681	27.66%	1.030%
83	19.491	2851	2860	2867	rBV	822030	1044116	47.68%	1.775%
84	19.697	2889	2895	2901	rBV	1837046	2189961	100.00%	3.723%
85	19.949	2933	2938	2945	rBV	86873	117608	5.37%	0.200%
86	20.396	3008	3014	3018	rBV	789223	877579	40.07%	1.492%
87	20.467	3018	3026	3030	rBV	939295	1164094	53.16%	1.979%
88	20.960	3105	3110	3120	rBV	213585	279263	12.75%	0.475%
89	21.172	3141	3146	3152	rBV2	1412284	1718385	78.47%	2.922%
90	21.236	3152	3157	3162	rVV2	937218	1580835	72.19%	2.688%
91	21.289	3162	3166	3171	rVB	954754	1067166	48.73%	1.814%
92	21.512	3200	3204	3211	rVB2	81732	108381	4.95%	0.184%
93	22.006	3284	3288	3290	rBV	43890	57046	2.60%	0.097%
94	22.041	3290	3294	3300	rVB	893630	1087672	49.67%	1.849%
95	22.817	3419	3426	3430	rBV	653631	1110482	50.71%	1.888%
96	22.864	3430	3434	3441	rVB	694117	1030844	47.07%	1.753%
97	23.398	3517	3525	3532	rVB	560140	1024639	46.79%	1.742%
98	23.492	3534	3541	3547	rBV	255584	468966	21.41%	0.797%
99	25.796	3918	3933	3947	rVB3	653791	2045676	93.41%	3.478%
100	26.489	4039	4051	4061	rBV2	315512	975927	44.56%	1.659%

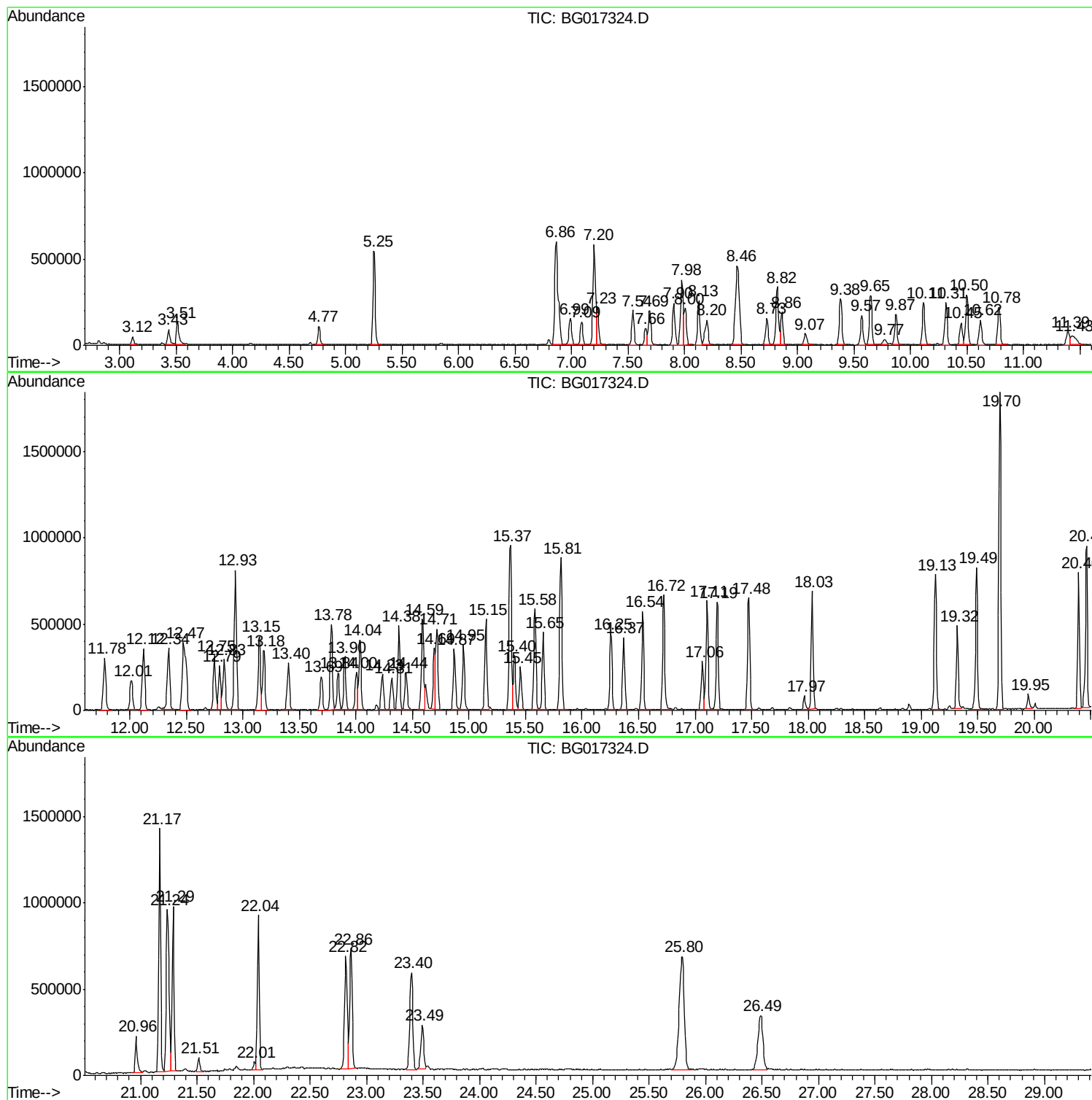
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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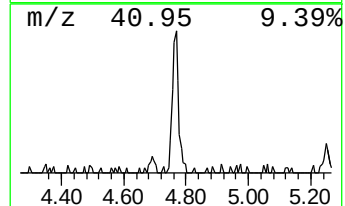
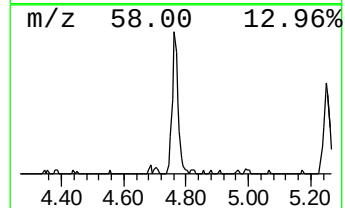
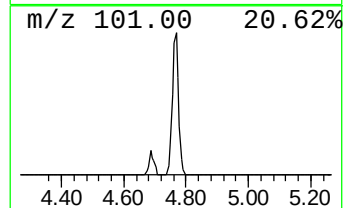
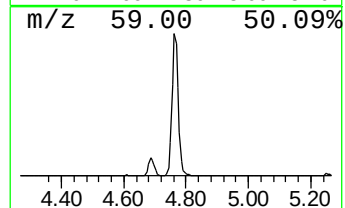
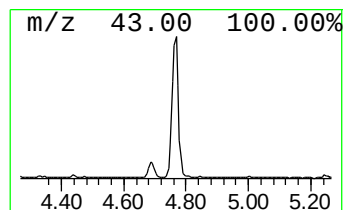
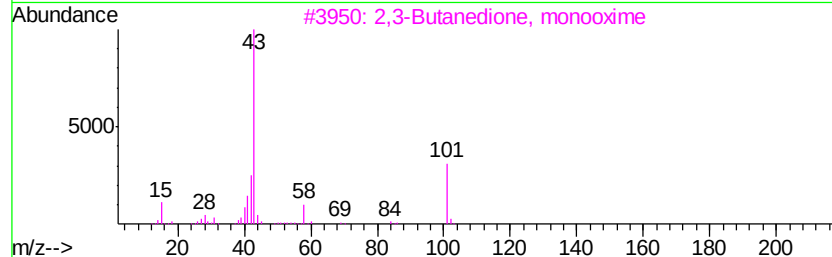
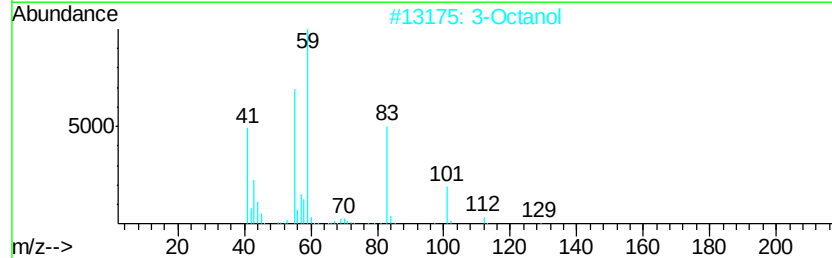
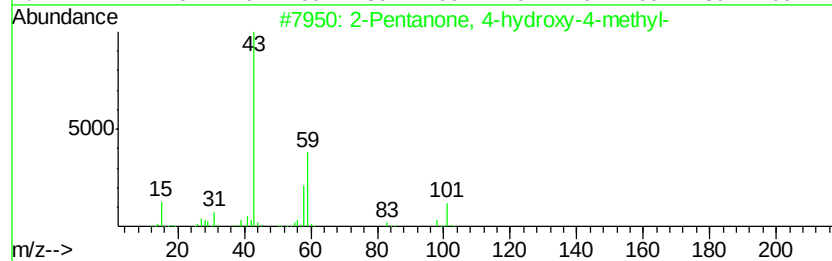
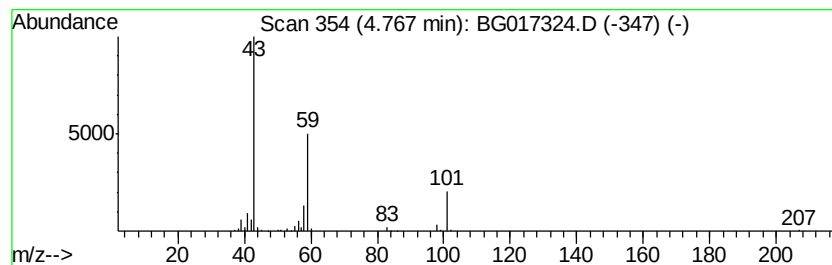
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.77	20.74 ng	156681	1,4-Dichlorobenzene-d4	7.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		3-Octanol	130	C8H18O	020296-29-1	33
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Propanenitrile, 3-ethoxy-	99	C5H9NO	002141-62-0	9
5		Dimethadione	129	C5H7NO3	000695-53-4	9



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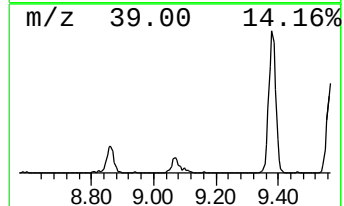
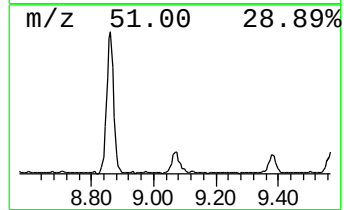
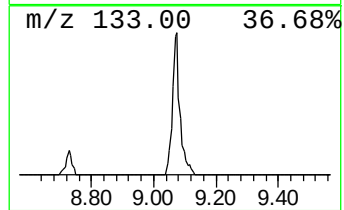
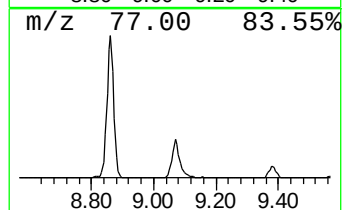
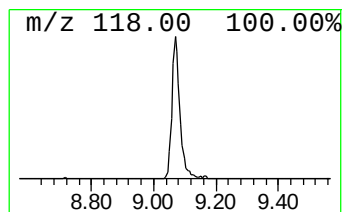
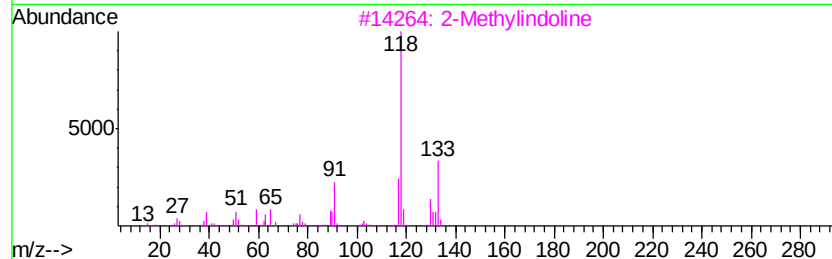
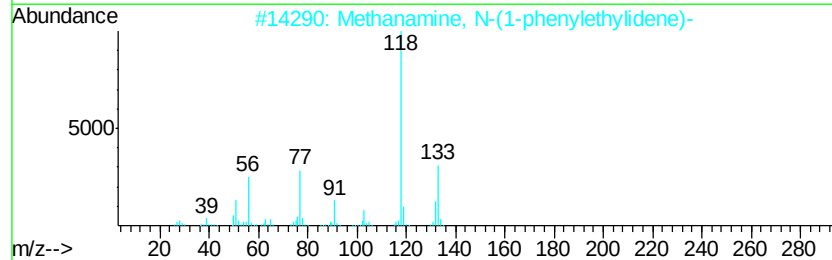
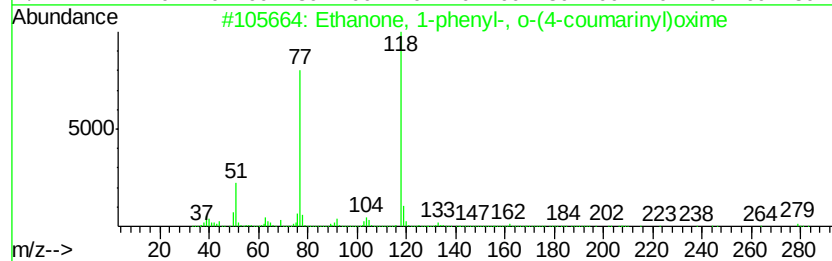
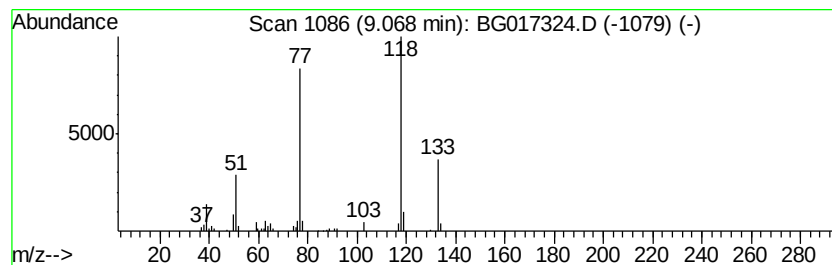
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Ethanone, 1-phenyl-, o-(4-c... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.07	10.99 ng	116456	Napthalene-d8	10.45
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Ethanone, 1-phenyl-, o-(4-coumar...	279 C17H13N03	034605-11-3	72
2	Methanamine, N-(1-phenylethylidene...	133 C9H11N	006907-71-7	40
3	2-Methylindoline	133 C9H11N	006872-06-6	38
4	Benzenecarboximidoyl bromide, N-...	197 C8H8BrN	041182-85-8	32
5	Benzonitrile, 2-amino-	118 C7H6N2	001885-29-6	27



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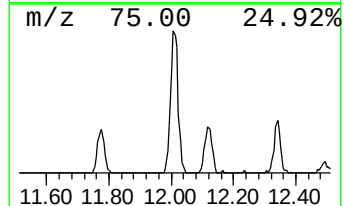
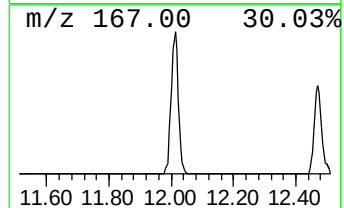
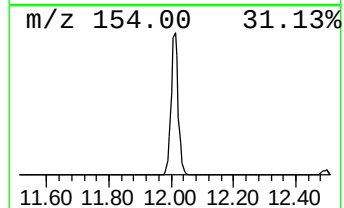
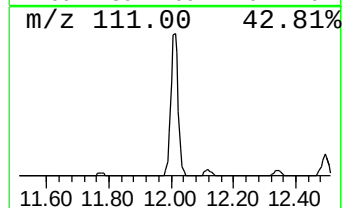
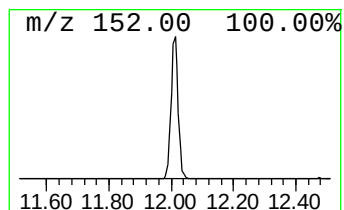
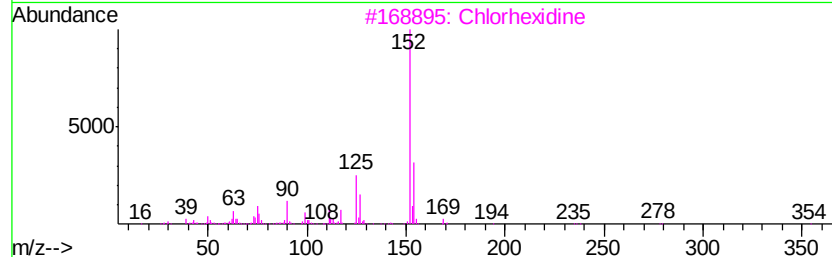
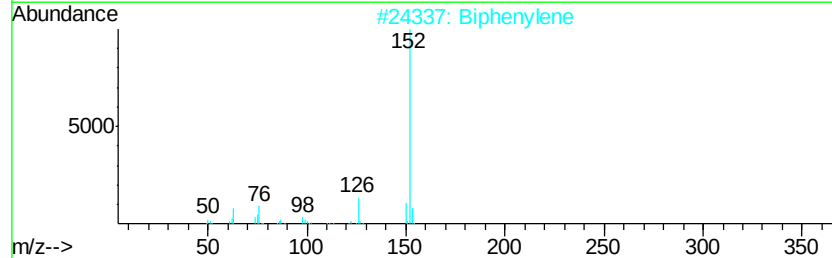
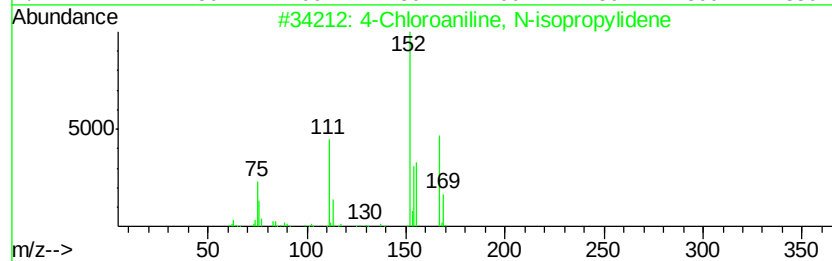
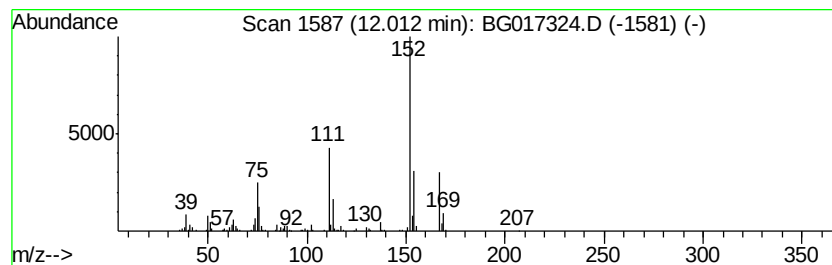
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 4-Chloroaniline, N-isopropy... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	26.29 ng	278443	Napthalene-d8	10.45

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4-Chloroaniline, N-isopropylidene	167	C9H10ClN	040938-43-0	91
2		Biphenylene	152	C12H8	000259-79-0	38
3		Chlorhexidine	504	C22H30Cl2N10	000055-56-1	37
4		2-Ethoxy-4-pyridinecarboxylic acid	167	C8H9NO3	091940-86-2	37
5		2-Chlorobenzimidazole	152	C7H5ClN2	004857-06-1	35



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG052815\
Data File : BG017324.D
Acq On : 28 May 2015 14:30
Operator : TP/IZ
Sample : G2387-09MS
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-18-D13MS

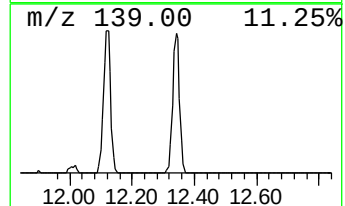
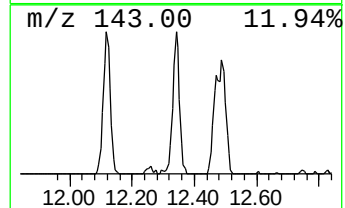
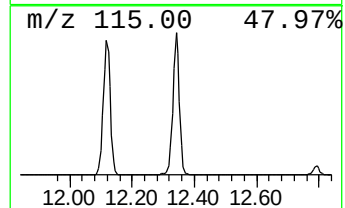
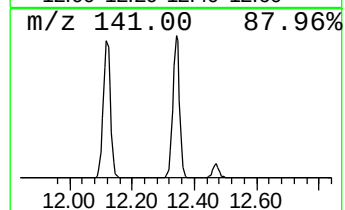
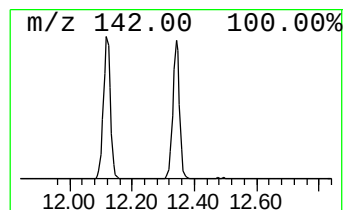
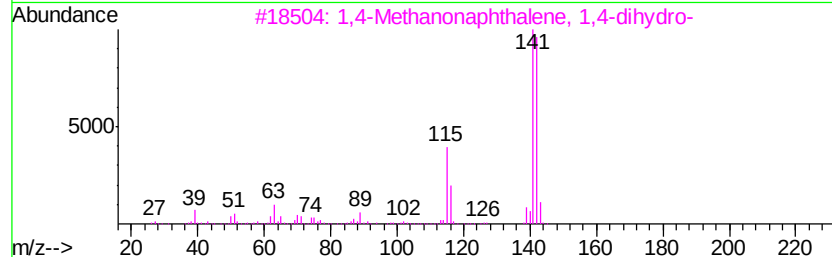
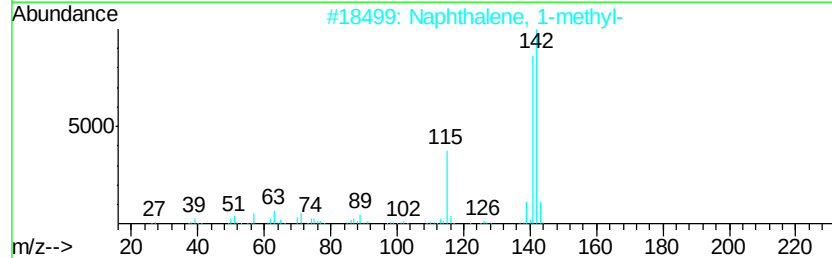
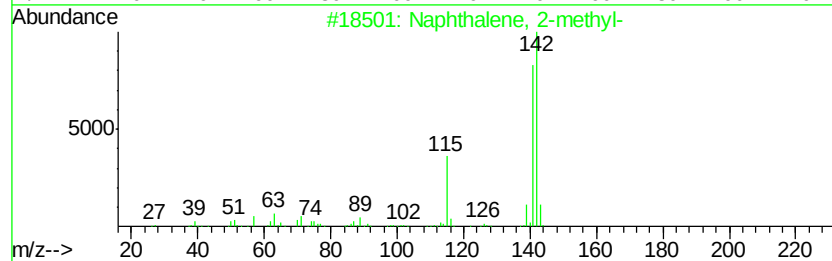
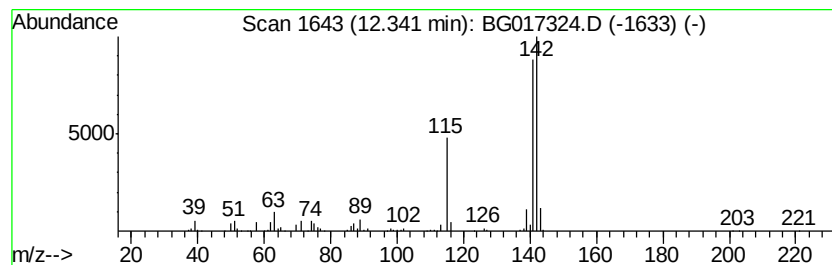
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Naphthalene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.34	53.47 ng	566326	Naphthalene-d8	10.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	97
3		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	59



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG052815\
Data File : BG017324.D
Acq On : 28 May 2015 14:30
Operator : TP/IZ
Sample : G2387-09MS
Misc :
ALS Vial : 5 Sample Multiplier: 1

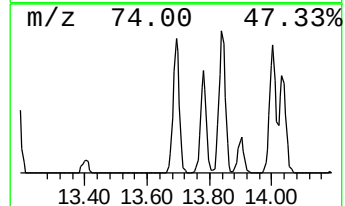
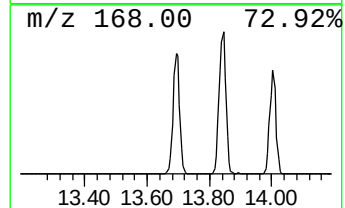
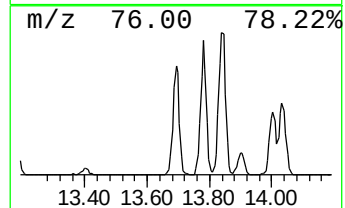
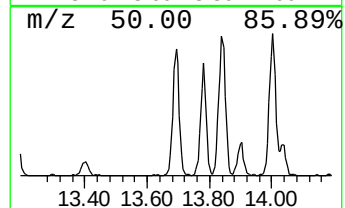
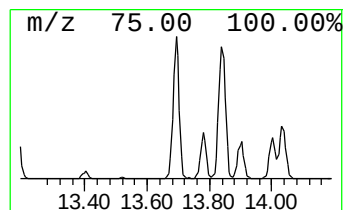
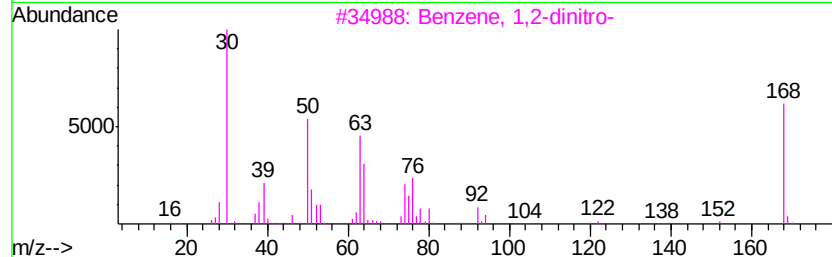
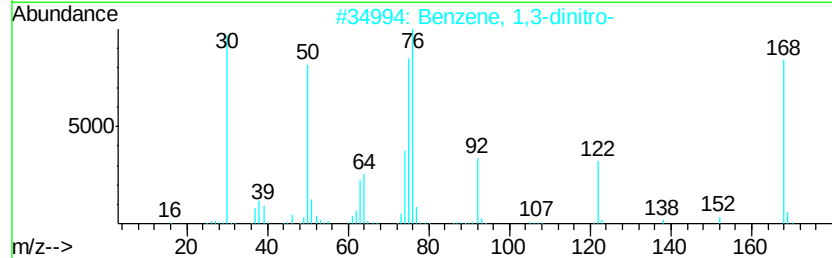
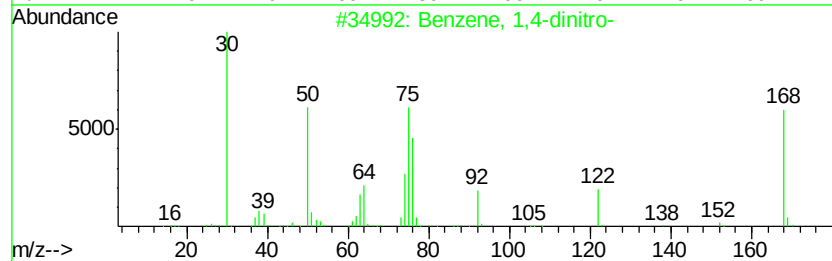
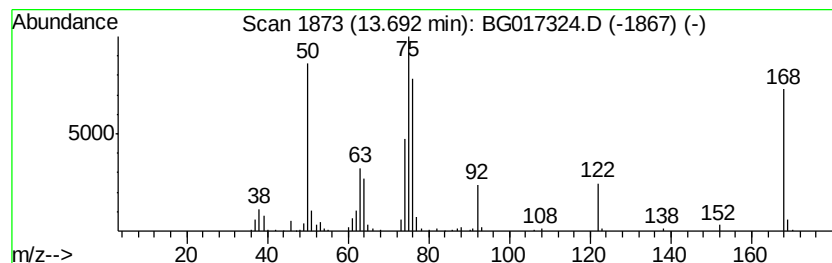
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ClientSampleId :
TP-18-D13MS

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzene, 1,4-dinitro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.69	19.19 ng	280208	Acenaphthene-d10	14.31
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,4-dinitro-	168 C6H4N2O4	000100-25-4 96
2		Benzene, 1,3-dinitro-	168 C6H4N2O4	000099-65-0 87
3		Benzene, 1,2-dinitro-	168 C6H4N2O4	000528-29-0 35
4		Bicyclo[4.2.0]octa-1,3,5-triene...	132 C8H4O2	006383-11-5 12
5		Silanamine, N-silyl-	77 H7NSi2	005702-11-4 9



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG052815\
Data File : BG017324.D
Acq On : 28 May 2015 14:30
Operator : TP/IZ
Sample : G2387-09MS
Misc :
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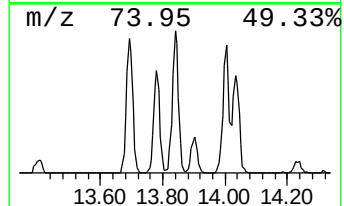
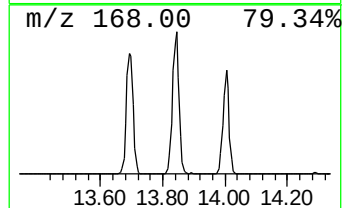
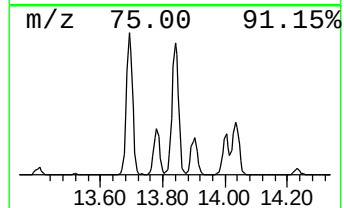
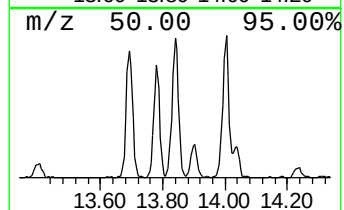
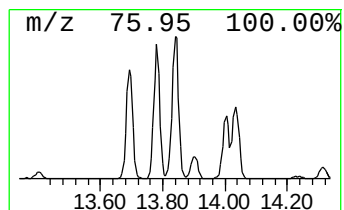
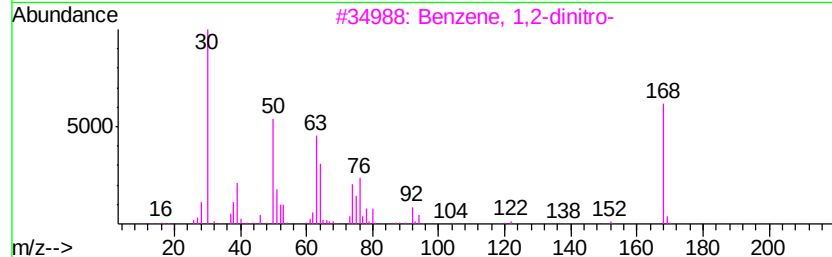
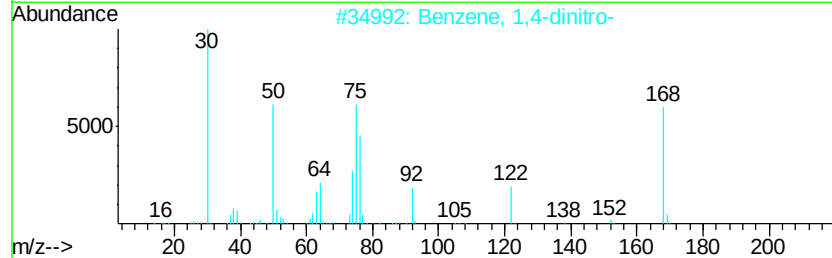
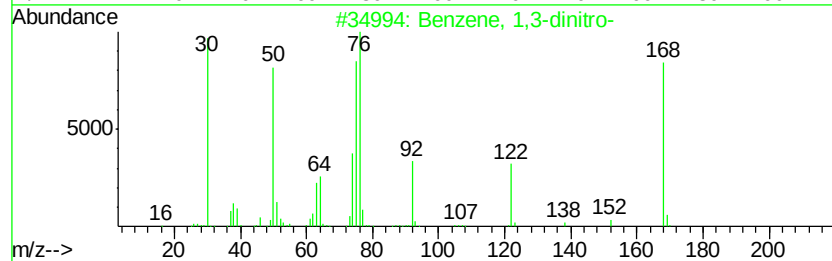
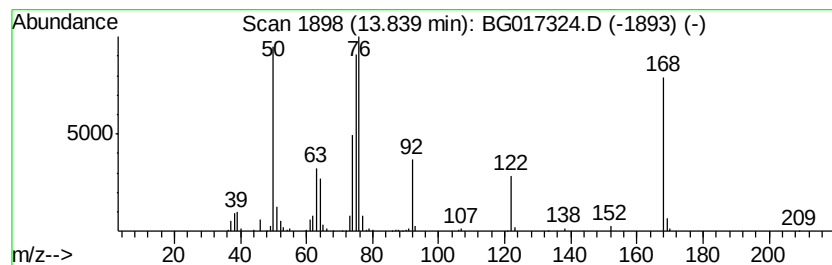
Instrument :
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ClientSampleId :
TP-18-D13MS

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, 1,3-dinitro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.84	22.26 ng	324955	Acenaphthene-d10	14.31
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,3-dinitro-	168 C6H4N2O4	000099-65-0 96
2		Benzene, 1,4-dinitro-	168 C6H4N2O4	000100-25-4 64
3		Benzene, 1,2-dinitro-	168 C6H4N2O4	000528-29-0 38
4		Benzoxazol-2-one, 3-(4-bromobenz...	351 C14H7BrClN03	1000269-38-7 13
5		Phthalic anhydride	148 C8H4O3	000085-44-9 9



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG052815\
Data File : BG017324.D
Acq On : 28 May 2015 14:30
Operator : TP/IZ
Sample : G2387-09MS
Misc :
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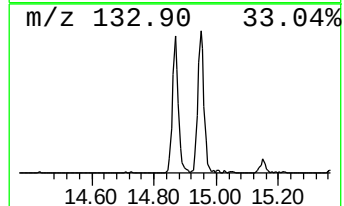
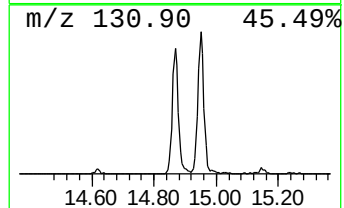
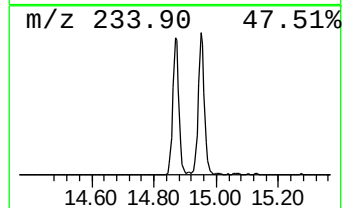
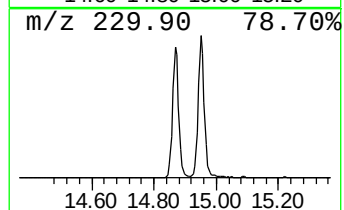
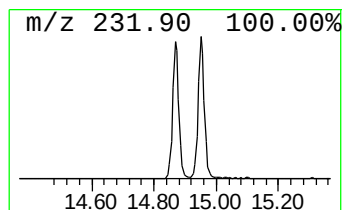
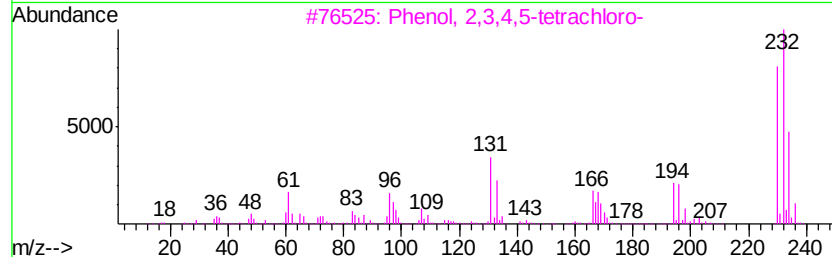
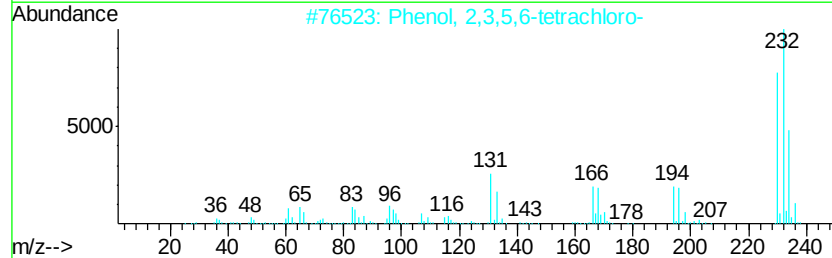
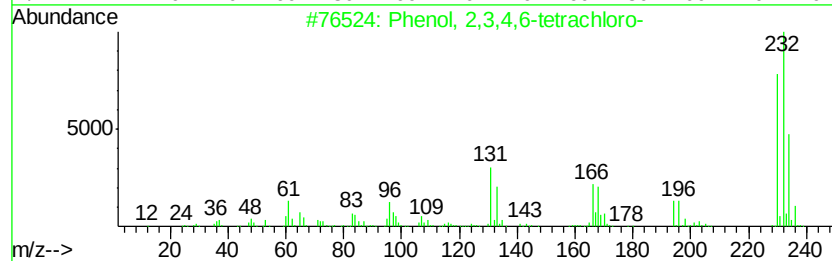
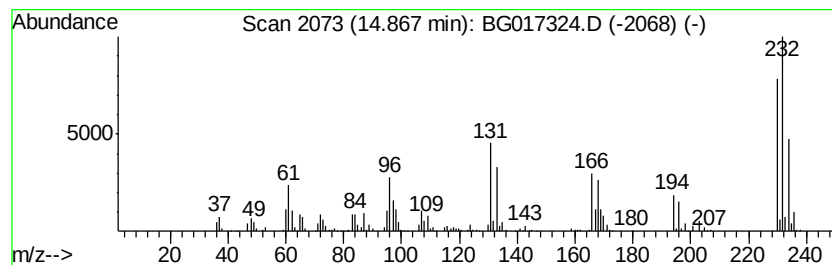
Instrument :
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ClientSampleId :
TP-18-D13MS

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenol, 2,3,4,6-tetrachloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	33.06 ng	482675	Acenaphthene-d10	14.31
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2 99
2	Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5 95
3	Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3 90
4	2,4,5,6-Tetrachloro-pyridin-3-yl...	230	C5H2Cl4N2	1000277-70-7 60
5	3-Bromo-7-methoxy pyrazolo[4,3-d...	228	C6H5BrN4O	068479-22-1 38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG052815\
Data File : BG017324.D
Acq On : 28 May 2015 14:30
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Sample : G2387-09MS
Misc :
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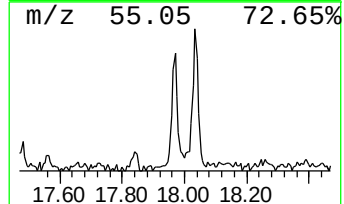
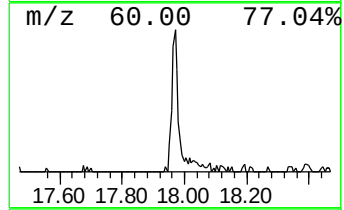
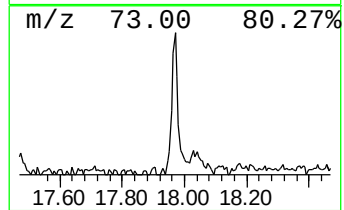
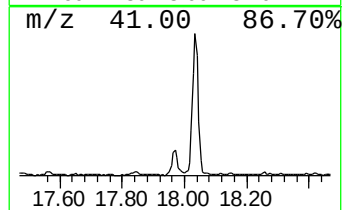
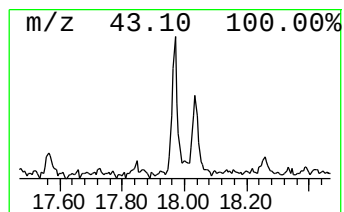
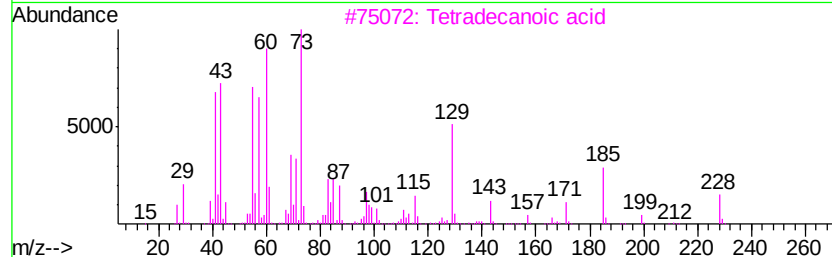
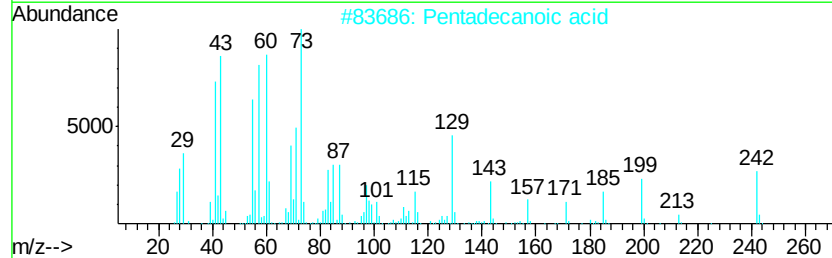
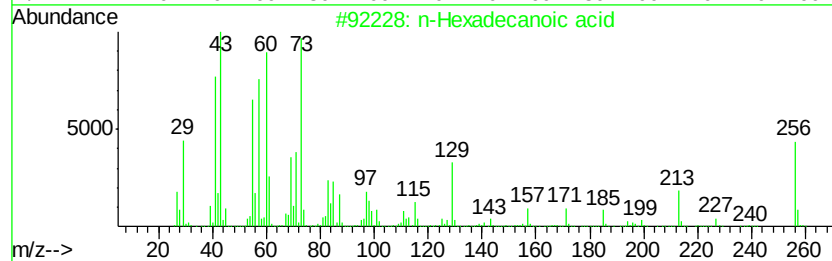
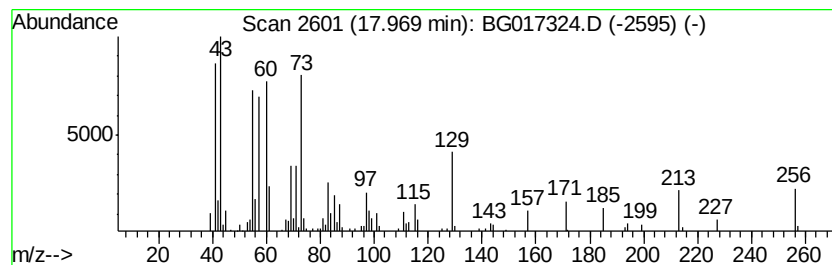
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.97	5.49 ng	109252	Phenanthrene-d10		17.06	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	53
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	53
4		Tridecanoic acid	214	C13H26O2	000638-53-9	47
5		n-Decanoic acid	172	C10H20O2	000334-48-5	46



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG052815\
Data File : BG017324.D
Acq On : 28 May 2015 14:30
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Sample : G2387-09MS
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TP-18-D13MS

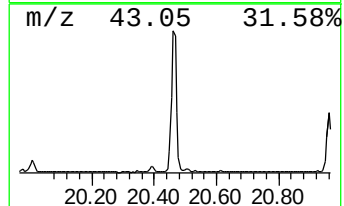
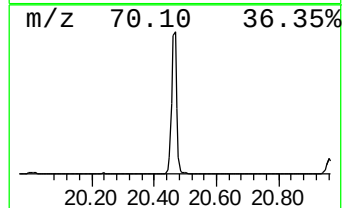
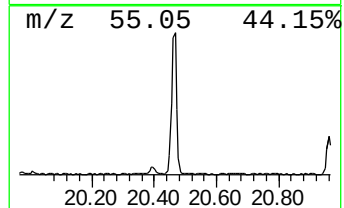
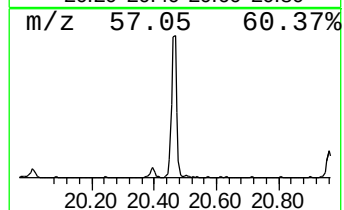
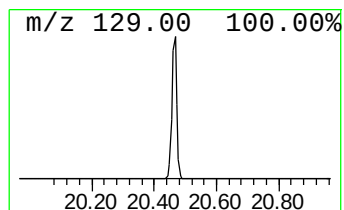
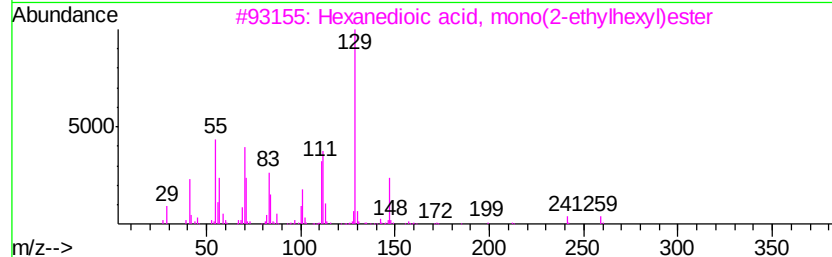
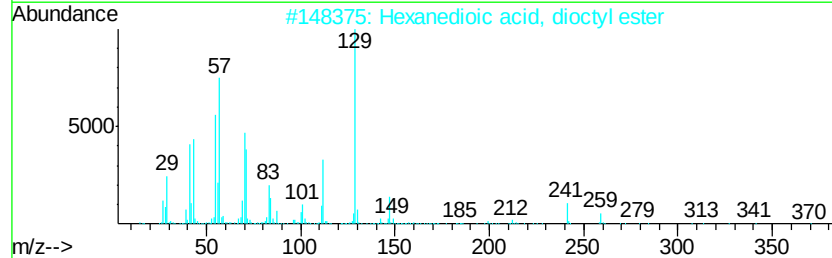
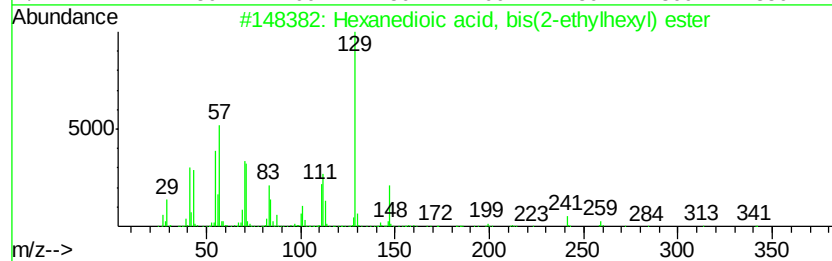
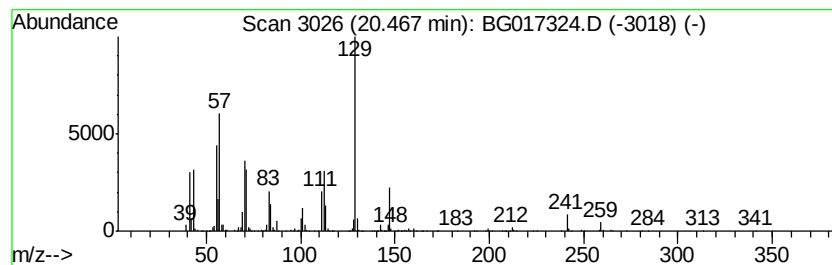
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexanedioic acid, bis(2-eth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.47	14.73 ng	1164090	Chrysene-d12	21.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	95
2		Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	93
3		Hexanedioic acid, mono(2-ethylhe...	258	C14H26O4	004337-65-9	58
4		Diisooctyl adipate	370	C22H42O4	001330-86-5	58
5		Adipic acid, di(oct-4-yl ester)	370	C22H42O4	1000160-80-1	47



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG052815\
Data File : BG017324.D
Acq On : 28 May 2015 14:30
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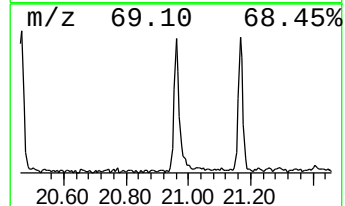
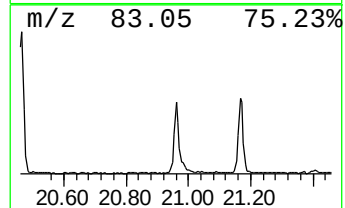
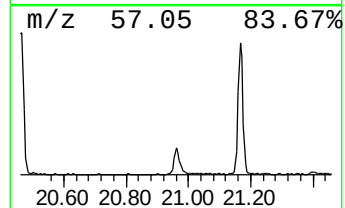
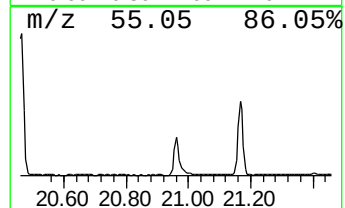
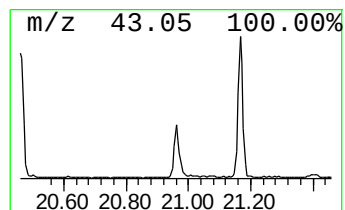
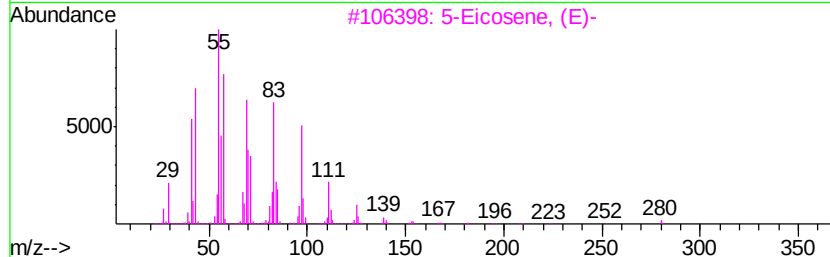
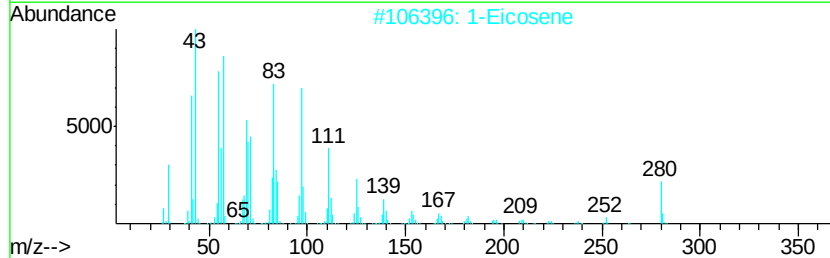
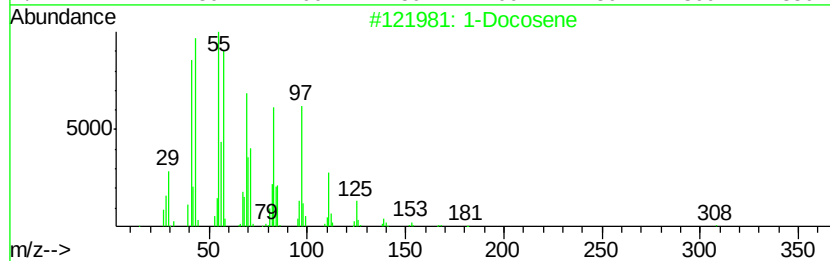
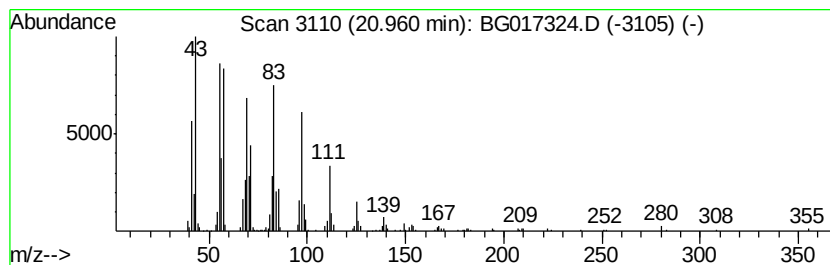
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BNA_G
ClientSampleId :
TP-18-D13MS

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1-Docosene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.96	3.53 ng	279263	Chrysene-d12	21.25
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Docosene		308 C22H44	001599-67-3 99
2	1-Eicosene		280 C20H40	003452-07-1 96
3	5-Eicosene, (E)-		280 C20H40	074685-30-6 96
4	1-Heneicosyl formate		340 C22H44O2	077899-03-7 95
5	Cyclohexadecane, 1,2-diethyl-		280 C20H40	1000155-85-3 94



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG052815\
Data File : BG017324.D
Acq On : 28 May 2015 14:30
Operator : TP/IZ
Sample : G2387-09MS
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TP-18-D13MS

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.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...	4.77	20.7	ng	156681	1	7.66	151091	20.0
Ethanone, 1-pheny...	9.07	11.0	ng	116456	2	10.45	211849	20.0
4-Chloroaniline, ...	12.01	26.3	ng	278443	2	10.45	211849	20.0
Naphthalene, 2-me...	12.34	53.5	ng	566326	2	10.45	211849	20.0
Benzene, 1,4-dini...	13.69	19.2	ng	280208	3	14.31	291974	20.0
Benzene, 1,3-dini...	13.84	22.3	ng	324955	3	14.31	291974	20.0
Phenol, 2,3,4,6-t...	14.87	33.1	ng	482675	3	14.31	291974	20.0
n-Hexadecanoic acid	17.97	5.5	ng	109252	4	17.06	398252	20.0
Hexanedioic acid, ...	20.47	14.7	ng	1164090	5	21.25	1580840	20.0
1-Docosene	20.96	3.5	ng	279263	5	21.25	1580840	20.0