

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122018\
 Data File : BG038770.D
 Acq On : 20 Dec 2018 20:12
 Operator : JU/SJ
 Sample : J6426-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P001-SS013-3036-01

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG121218.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	4.217	182	192	213	rBV	6039912	11200679	50.77%	11.925%
2	5.544	411	418	426	rBV	2023436	3766327	17.07%	4.010%
3	5.627	426	432	435	rVV	392377	691972	3.14%	0.737%
4	5.703	435	445	458	rVB	8303712	22062580	100.00%	23.490%
5	6.061	496	506	513	rVB	3929243	7089112	32.13%	7.548%
6	7.119	679	686	692	rBV	3010028	4965936	22.51%	5.287%
7	7.213	692	702	707	rVV2	415744	1072746	4.86%	1.142%
8	7.260	707	710	717	rVB	165607	267311	1.21%	0.285%
9	7.430	731	739	749	rBV	114022	252426	1.14%	0.269%
10	7.589	759	766	771	rBV	411370	727542	3.30%	0.775%
11	7.642	771	775	779	rVV	278014	420827	1.91%	0.448%
12	7.689	779	783	790	rVB	448054	766824	3.48%	0.816%
13	8.065	840	847	857	rVB5	125776	349102	1.58%	0.372%
14	8.171	857	865	875	rBV3	202544	503111	2.28%	0.536%
15	8.382	894	901	907	rBV	302484	551754	2.50%	0.587%
16	8.594	932	937	945	rVB2	119603	265966	1.21%	0.283%
17	8.682	945	952	959	rBV3	231737	435091	1.97%	0.463%
18	9.146	1026	1031	1039	rVB3	145809	278611	1.26%	0.297%
19	9.246	1039	1048	1055	rBV2	698295	1408337	6.38%	1.499%
20	9.557	1094	1101	1113	rVB	1712129	2764063	12.53%	2.943%
21	9.969	1153	1171	1177	rBV6	84245	313020	1.42%	0.333%
22	10.251	1214	1219	1226	rBV2	187543	346610	1.57%	0.369%
23	10.732	1291	1301	1306	rBV	227283	446609	2.02%	0.475%
24	10.803	1306	1313	1321	rVV	697236	1259510	5.71%	1.341%
25	10.873	1321	1325	1329	rVV3	207993	408789	1.85%	0.435%
26	10.926	1329	1334	1339	rVV	448091	862821	3.91%	0.919%
27	10.985	1339	1344	1350	rVB2	250357	451033	2.04%	0.480%
28	11.549	1432	1440	1445	rBV6	122514	292877	1.33%	0.312%
29	11.743	1469	1473	1483	rVB3	126124	295017	1.34%	0.314%
30	11.849	1484	1491	1495	rBV	302491	502399	2.28%	0.535%
31	12.049	1519	1525	1531	rBV	644810	1052886	4.77%	1.121%
32	12.248	1552	1559	1567	rBV	978516	1557538	7.06%	1.658%
33	12.524	1601	1606	1612	rBV	445889	754257	3.42%	0.803%
34	12.748	1635	1644	1651	rVB	234215	470816	2.13%	0.501%

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35	12.871	1660	1665	1670	rBV4	121329	237201	1.08%	0.253%
36	13.053	1689	1696	1702	rVB3	226626	405184	1.84%	0.431%
37	13.188	1714	1719	1725	rBV	293224	442604	2.01%	0.471%
38	13.312	1735	1740	1749	rVB	624833	1047526	4.75%	1.115%
39	13.482	1762	1769	1775	rBV2	1130085	1891566	8.57%	2.014%
40	13.723	1805	1810	1819	rVB2	145814	365054	1.65%	0.389%
41	13.870	1827	1835	1841	rVB4	292976	720403	3.27%	0.767%
42	14.017	1853	1860	1865	rBV2	444412	1074721	4.87%	1.144%
43	14.064	1865	1868	1872	rVV	236675	409631	1.86%	0.436%
44	14.134	1876	1880	1884	rVV2	355069	553782	2.51%	0.590%
45	14.175	1884	1887	1894	rVB3	133942	233836	1.06%	0.249%
46	14.252	1895	1900	1904	rBV	234735	355649	1.61%	0.379%
47	14.416	1923	1928	1934	rBV3	148191	255421	1.16%	0.272%
48	14.551	1946	1951	1959	rVB	1079514	1443302	6.54%	1.537%
49	14.698	1972	1976	1981	rVB3	146843	278998	1.26%	0.297%
50	14.892	2004	2009	2018	rVB3	186947	418055	1.89%	0.445%
51	15.080	2037	2041	2048	rVB3	236796	466637	2.12%	0.497%
52	15.163	2049	2055	2063	rBV2	360573	655326	2.97%	0.698%
53	15.304	2073	2079	2084	rBV	189936	378172	1.71%	0.403%
54	15.497	2101	2112	2117	rBV	1025871	1750159	7.93%	1.863%
55	15.897	2172	2180	2184	rBV2	361349	717913	3.25%	0.764%
56	16.050	2203	2206	2214	rVB5	175978	258517	1.17%	0.275%
57	16.185	2224	2229	2235	rBV3	548368	899689	4.08%	0.958%
58	16.361	2254	2259	2276	rVB	986673	2300227	10.43%	2.449%
59	17.154	2389	2394	2398	rBV	674980	887295	4.02%	0.945%
60	17.201	2398	2402	2407	rVB2	367183	536267	2.43%	0.571%
61	17.442	2438	2443	2447	rBV	240401	411507	1.87%	0.438%
62	17.483	2447	2450	2455	rVB	235933	336898	1.53%	0.359%
63	17.577	2461	2466	2470	rBV	357516	482636	2.19%	0.514%
64	17.889	2515	2519	2529	rVB	612087	857778	3.89%	0.913%
65	18.582	2633	2637	2642	rVB	426042	490740	2.22%	0.522%
66	19.211	2740	2744	2751	rBV2	352275	516828	2.34%	0.550%
67	19.781	2838	2841	2846	rVB2	183842	227682	1.03%	0.242%
68	20.045	2881	2886	2890	rVB	1112971	1441183	6.53%	1.534%
69	20.315	2928	2932	2937	rBV2	223644	321090	1.46%	0.342%
70	21.725	3167	3172	3180	rVB	317622	607557	2.75%	0.647%
71	25.027	3728	3734	3744	rVB2	138516	393614	1.78%	0.419%

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

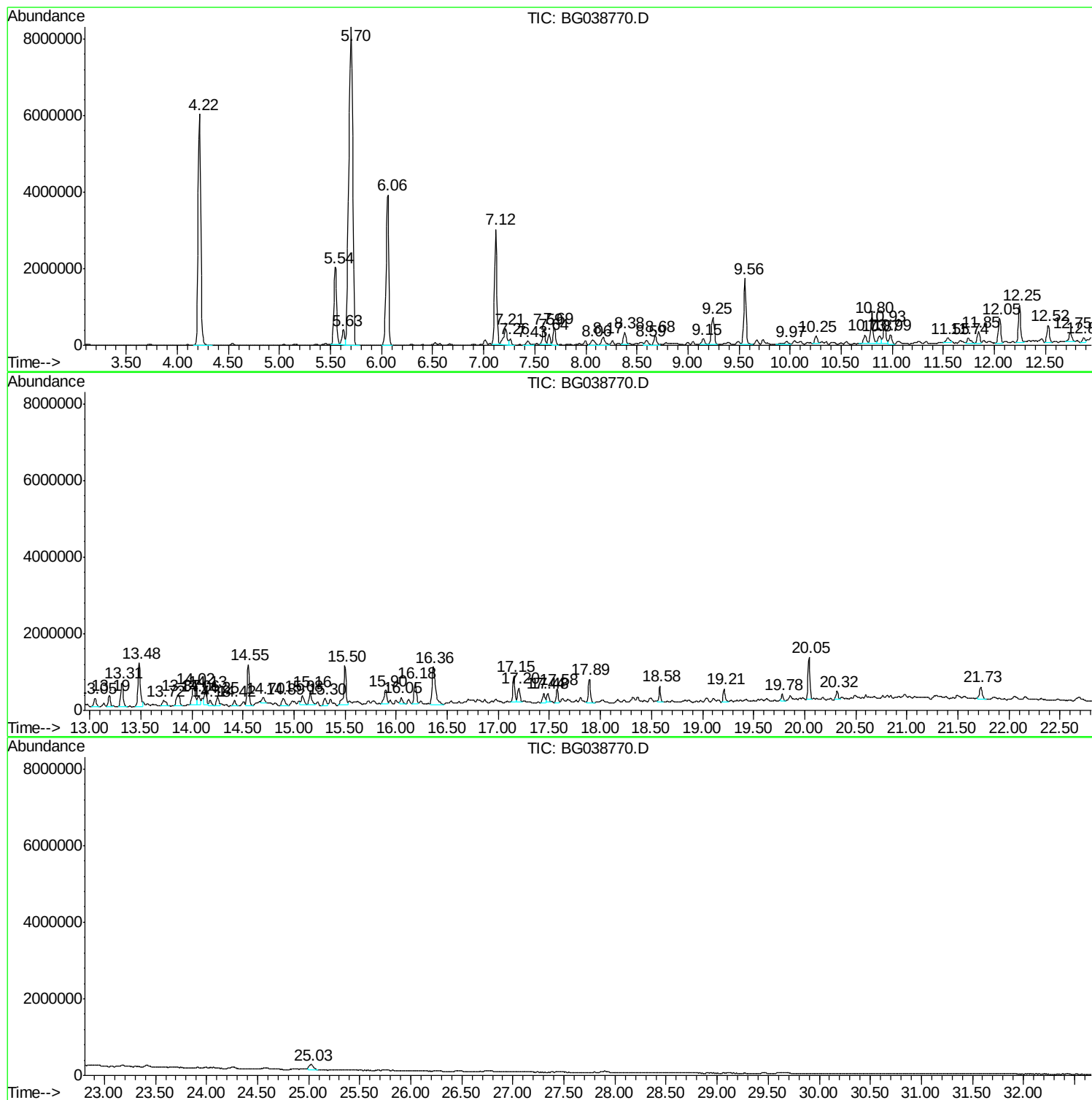
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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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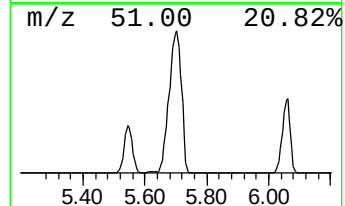
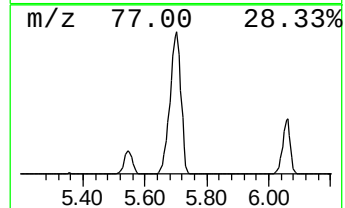
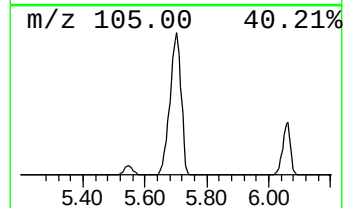
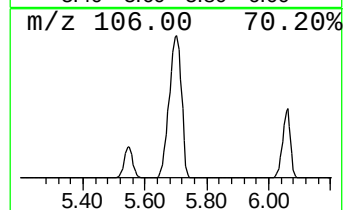
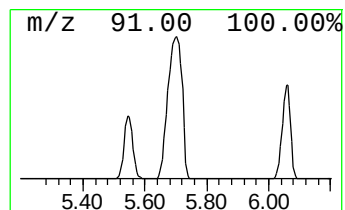
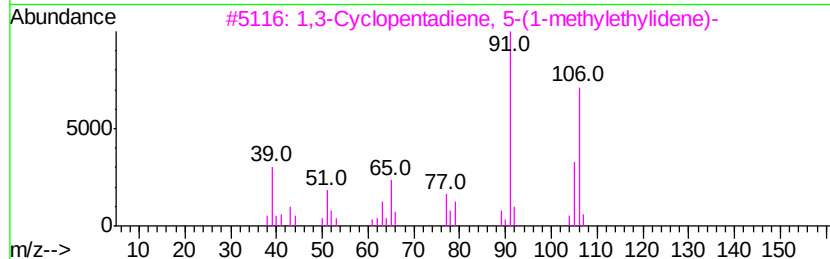
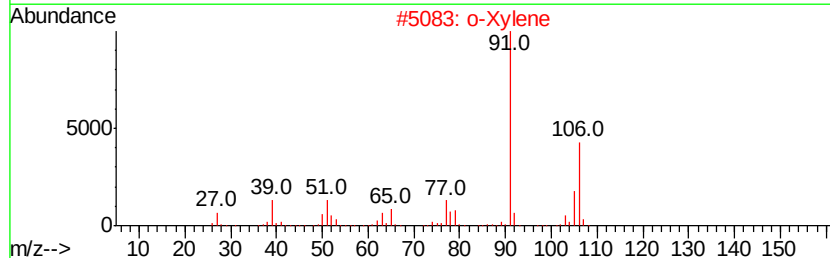
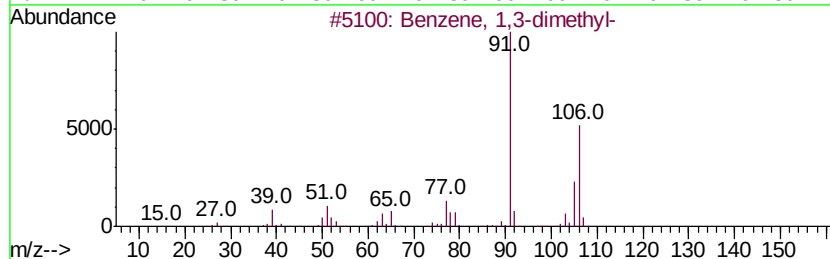
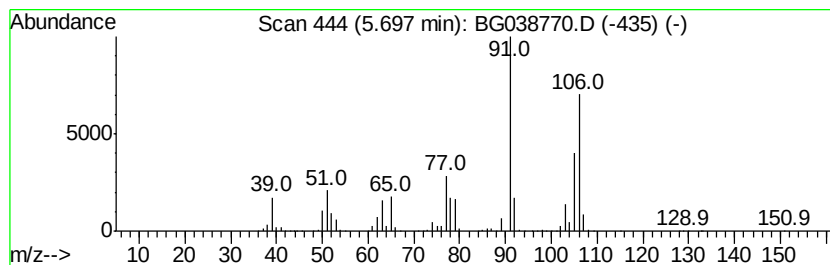
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.70	1263.96 ng	22062600	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	91
2		o-Xylene	106	C8H10	000095-47-6	90
3		1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	83
4		p-Xylene	106	C8H10	000106-42-3	81
5		Ethylbenzene	106	C8H10	000100-41-4	64



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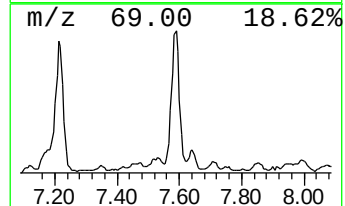
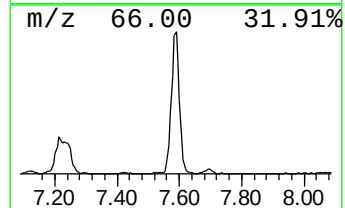
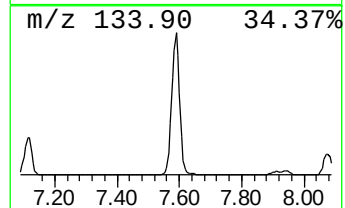
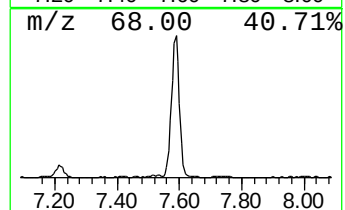
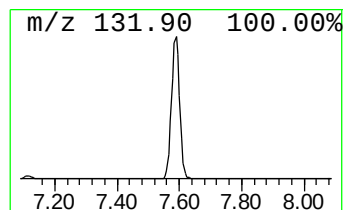
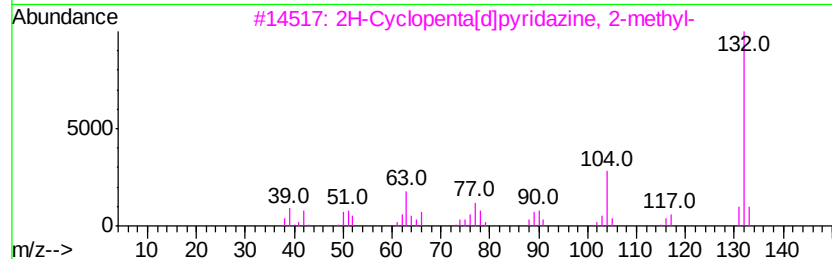
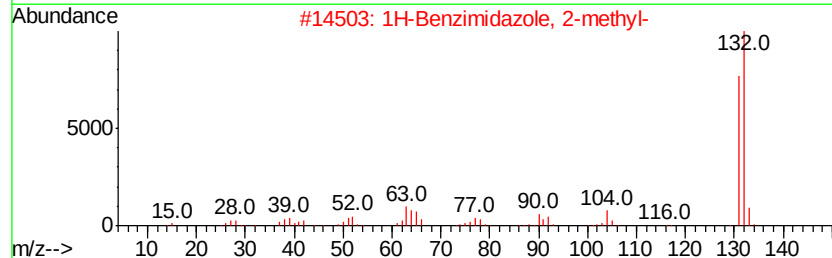
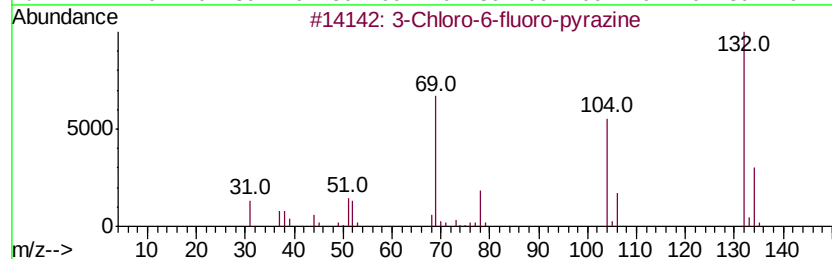
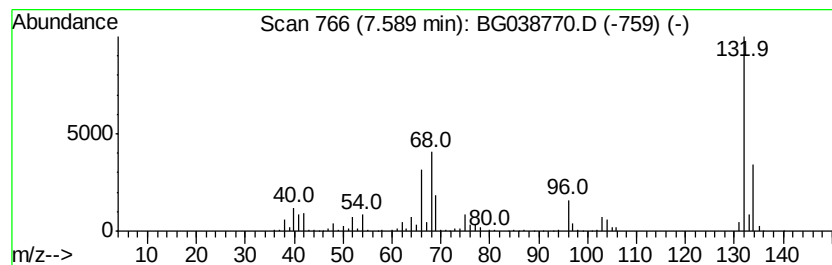
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
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TIC Library : C:\Database\NIST11.L
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Peak Number 6 unknown7.59 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	41.68 ng	727542	1,4-Dichlorobenzene-d4	8.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	25
3			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	22
4			5-Aminoindole	132	C8H8N2	005192-03-0	22
5			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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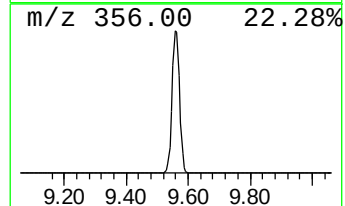
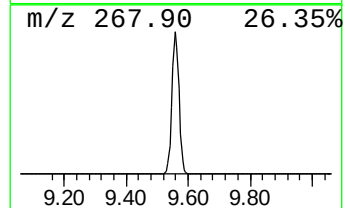
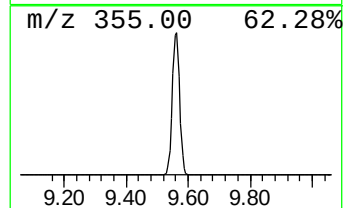
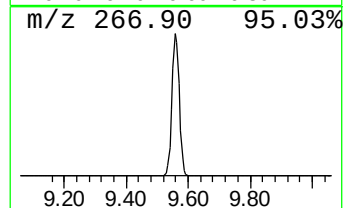
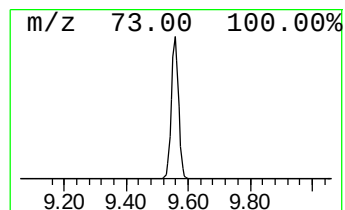
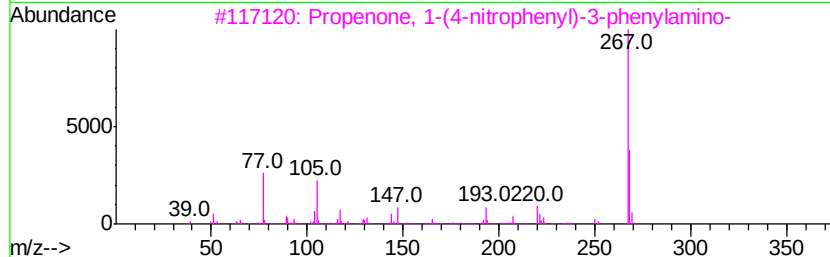
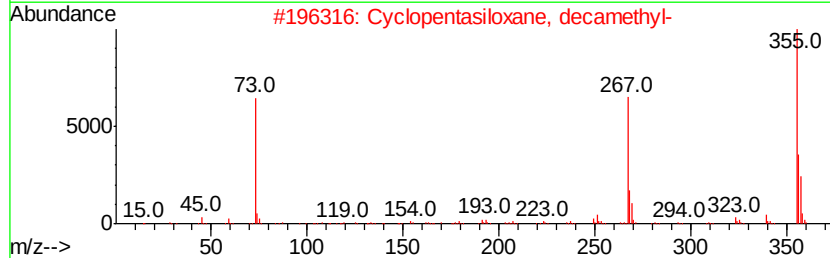
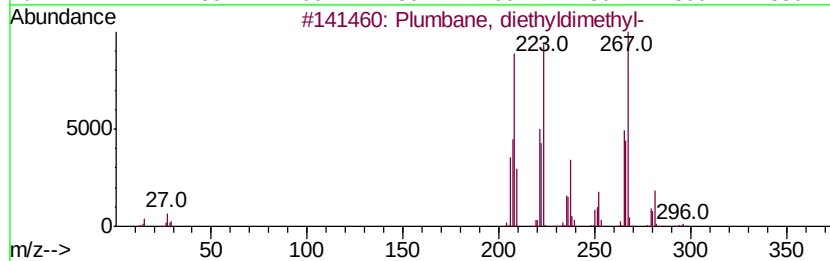
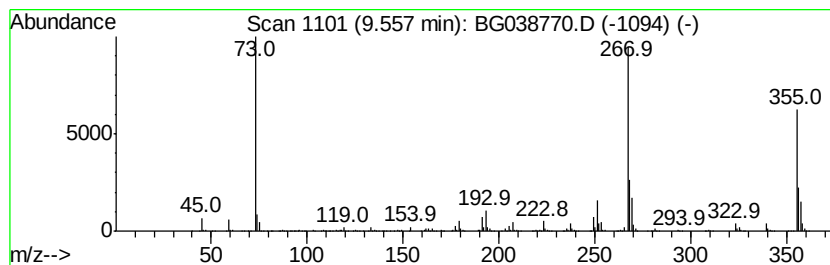
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TIC Library : C:\Database\NIST11.L
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Peak Number 8 Plumbane, diethyldimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.56	135.23 ng	2764060	Naphthalene-d8	10.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Plumbane, diethyldimethyl-	296	C6H16Pb	001762-27-2	94
2		Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	91
3		Propenone, 1-(4-nitrophenyl)-3-phenylamino-	268	C15H12N2O3	1000302-96-9	38
4		Benzaldehyde, 2,4-bis(trimethylsilyl)-	282	C13H22O3Si2	033617-38-8	37
5		Silane, [(fluoromethylphenyl)silyl]-	370	C17H35FSi4	072190-81-9	35



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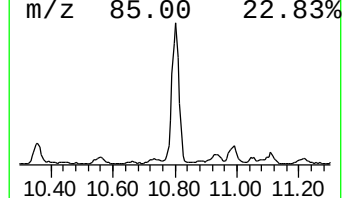
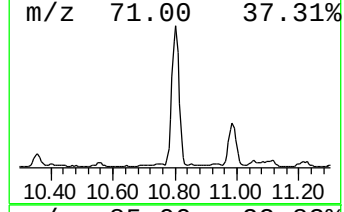
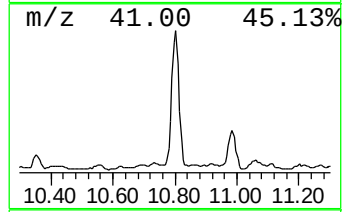
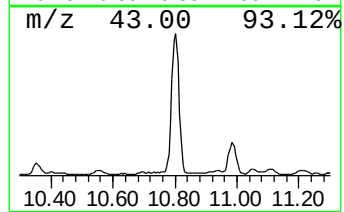
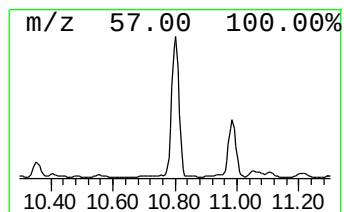
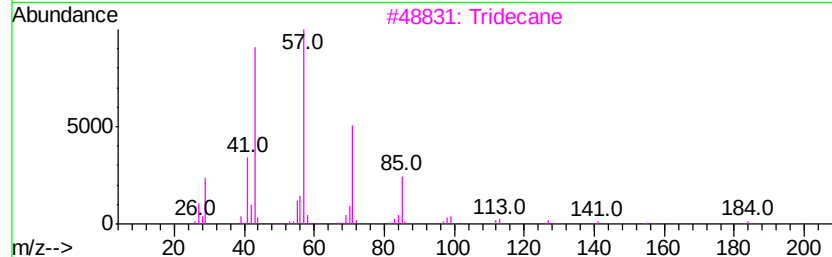
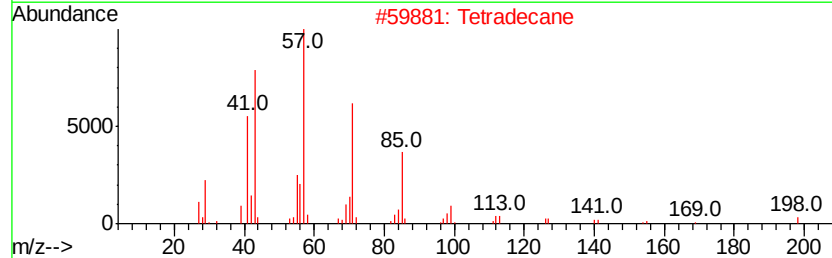
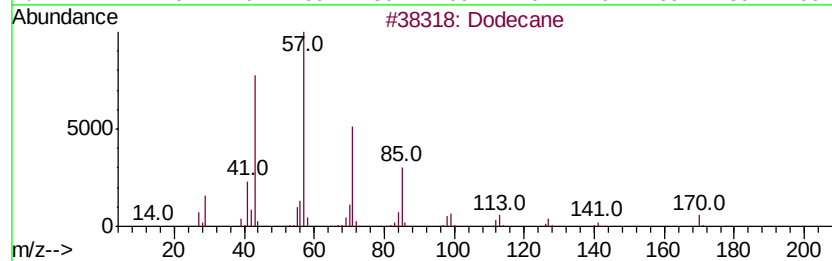
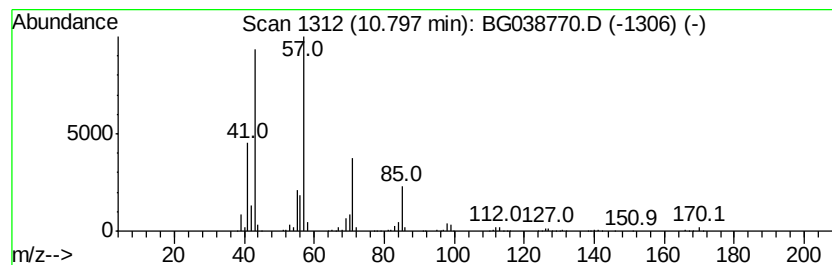
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Dodecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.80	61.62 ng	1259510	Naphthalene-d8	10.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane	170	C12H26	000112-40-3	90
2		Tetradecane	198	C14H30	000629-59-4	90
3		Tridecane	184	C13H28	000629-50-5	86
4		Undecane	156	C11H24	001120-21-4	83
5		Hexadecane	226	C16H34	000544-76-3	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122018\
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Acq On : 20 Dec 2018 20:12
Operator : JU/SJ
Sample : J6426-08
Misc :
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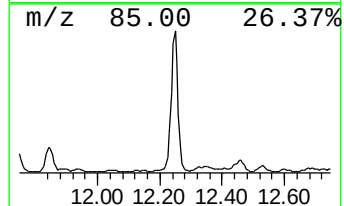
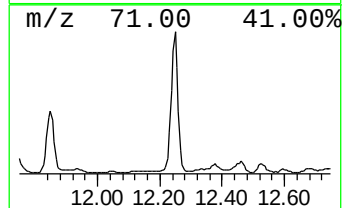
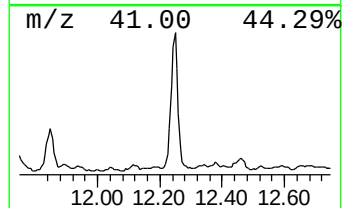
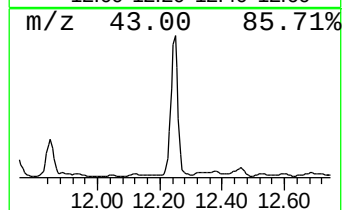
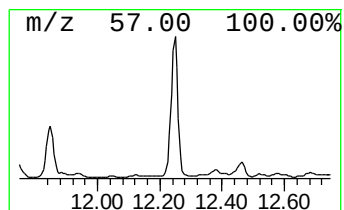
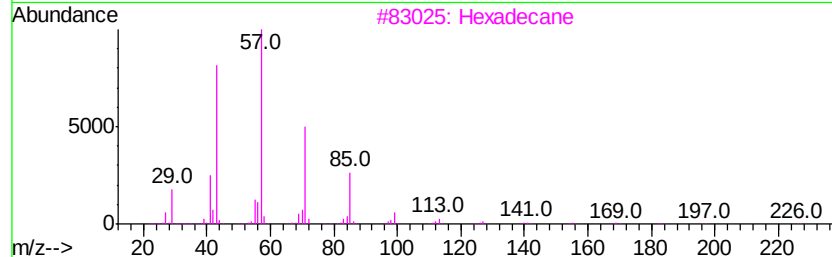
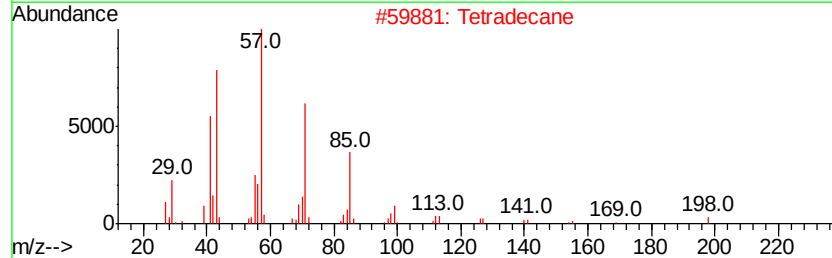
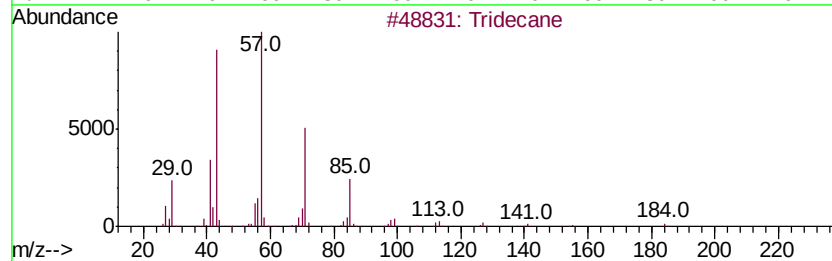
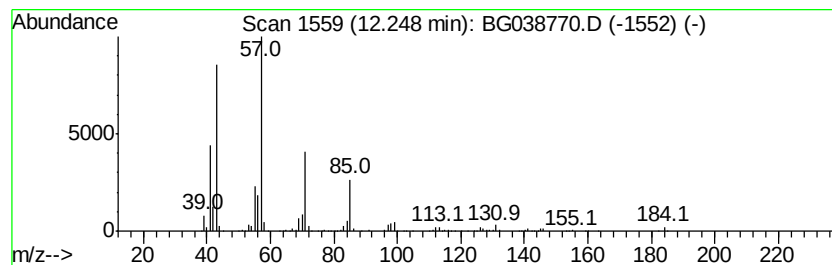
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tridecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.25	76.20 ng	1557540	Naphthalene-d8	10.87	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	95
2	Tetradecane	198	C14H30	000629-59-4	90
3	Hexadecane	226	C16H34	000544-76-3	90
4	Dodecane	170	C12H26	000112-40-3	83
5	Pentadecane	212	C15H32	000629-62-9	83



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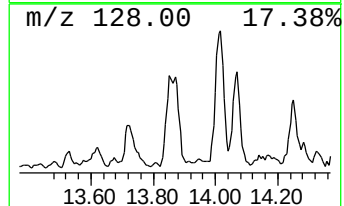
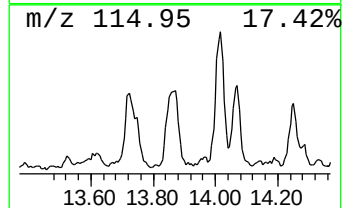
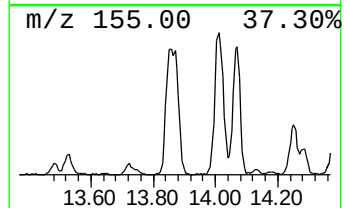
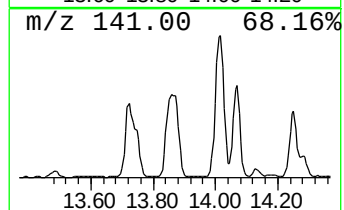
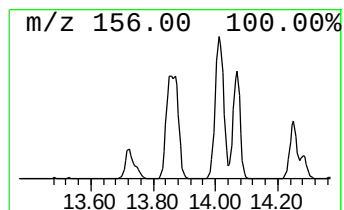
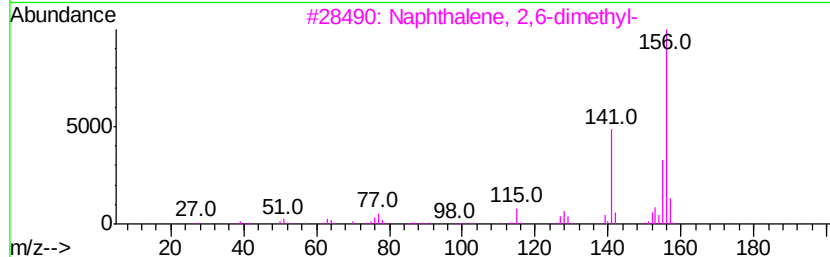
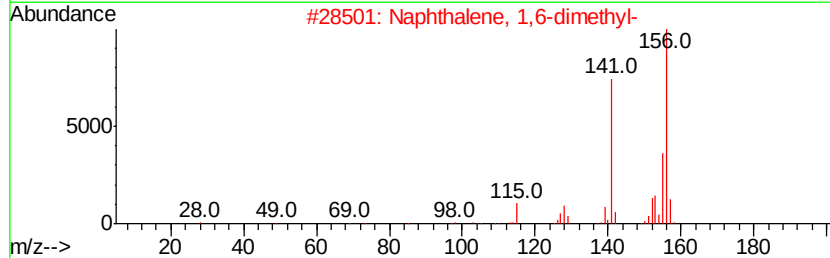
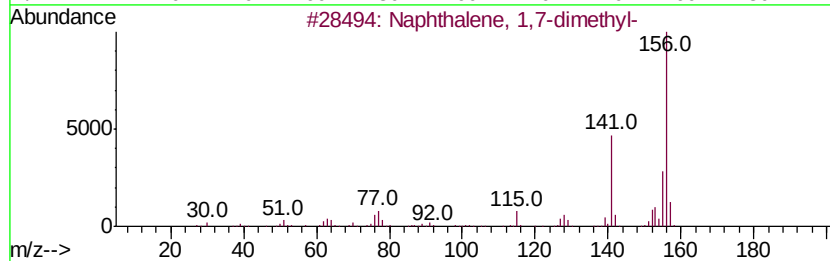
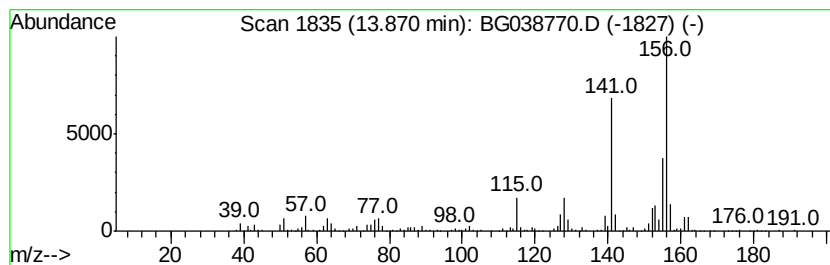
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Naphthalene, 1,7-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.		
13.87	51.64 ng	720403	Acenaphthene-d10		14.69		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	95
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
5			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122018\
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Operator : JU/SJ
Sample : J6426-08
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Instrument :
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ClientSampleId :
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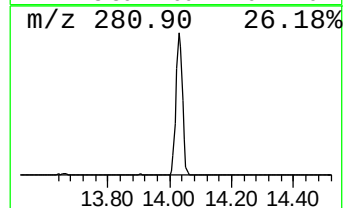
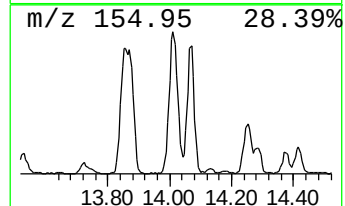
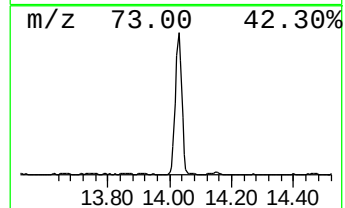
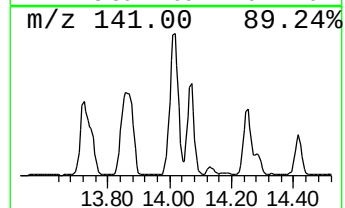
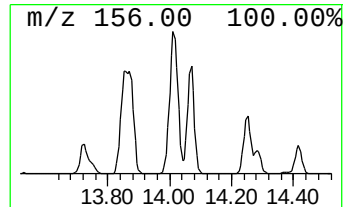
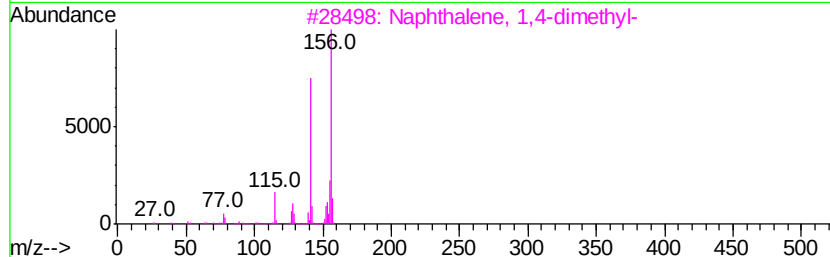
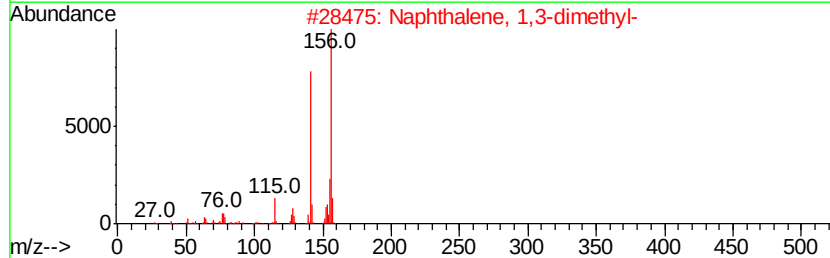
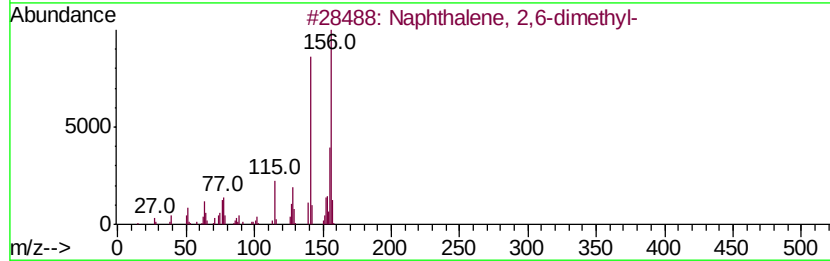
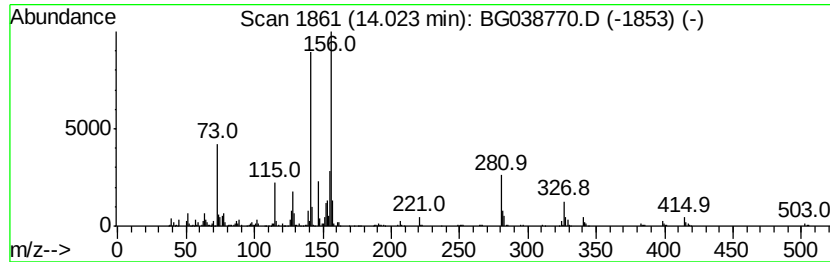
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Naphthalene, 2,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.02	77.04 ng	1074720	Acenaphthene-d10	14.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	95
2		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
3		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
4		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
5		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122018\
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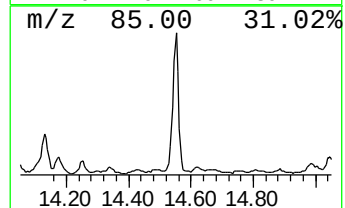
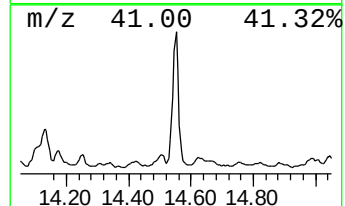
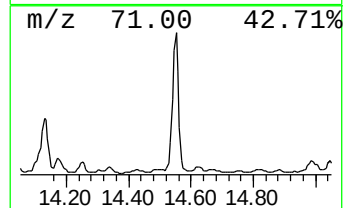
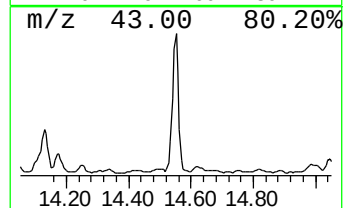
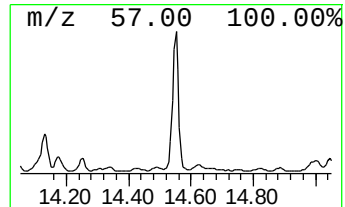
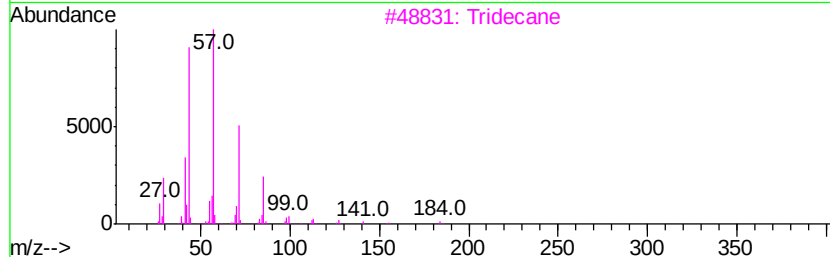
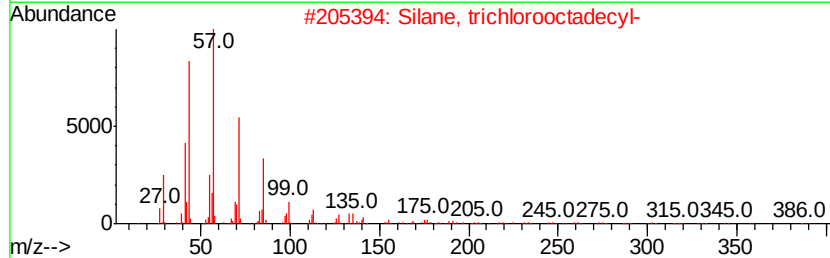
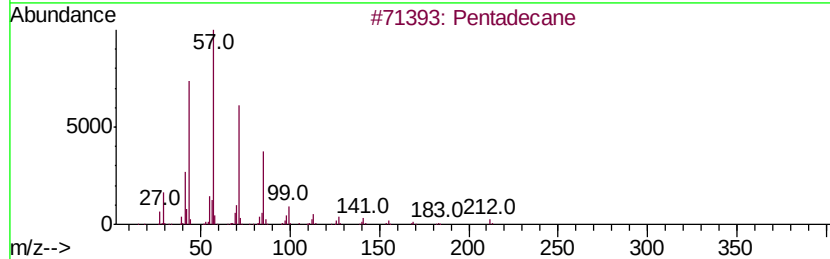
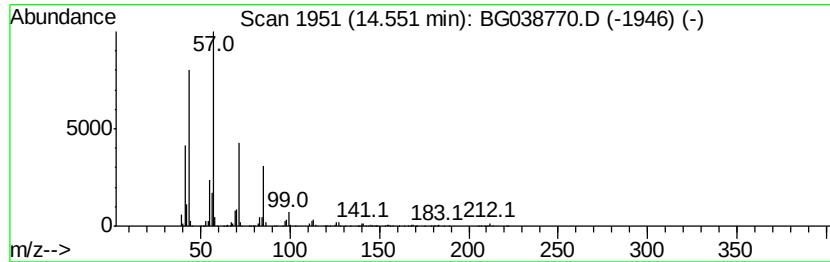
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Pentadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.55	103.46 ng	1443300	Acenaphthene-d10	14.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane	212	C15H32	000629-62-9	93
2		Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	90
3		Tridecane	184	C13H28	000629-50-5	90
4		Hexadecane	226	C16H34	000544-76-3	90
5		Undecane	156	C11H24	001120-21-4	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122018\
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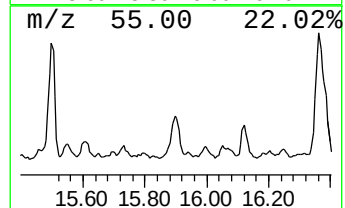
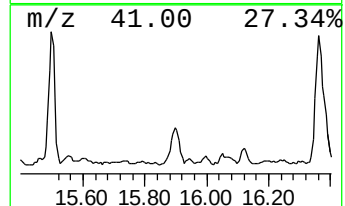
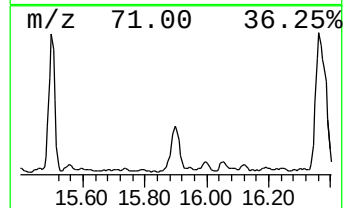
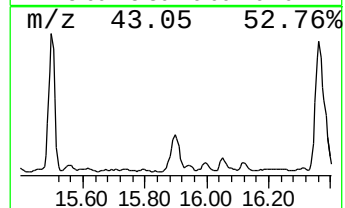
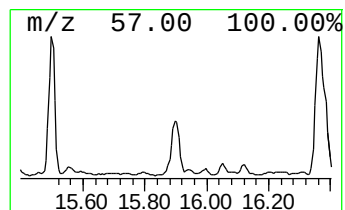
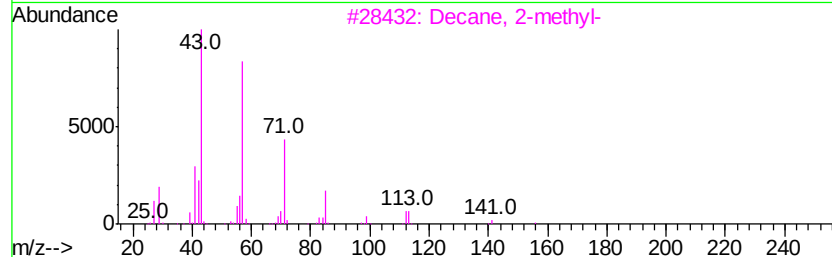
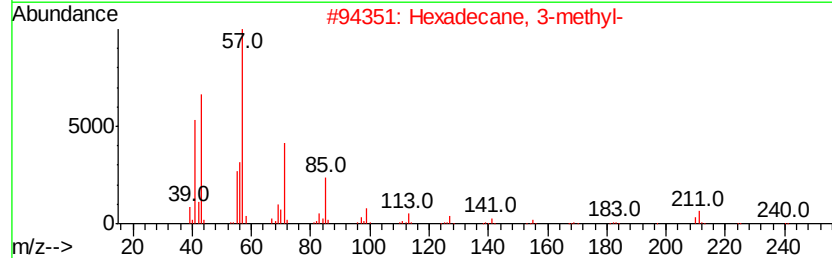
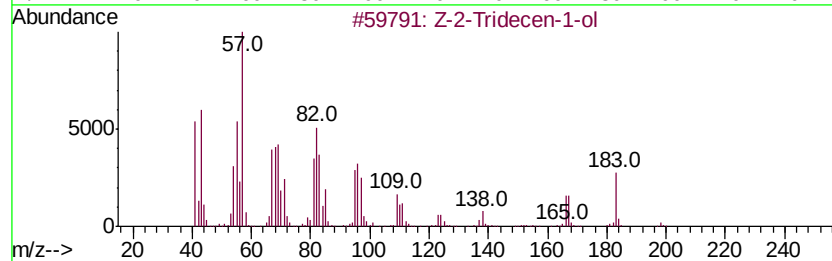
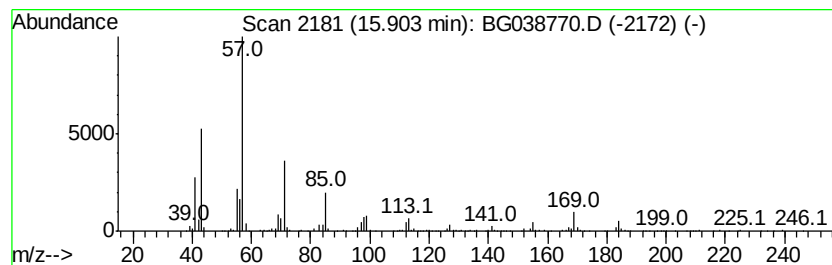
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Z-2-Tridecen-1-ol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.90	51.46 ng	717913	Acenaphthene-d10		14.69	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Z-2-Tridecen-1-ol		198	C13H26O	1000130-75-8	83
2	Hexadecane, 3-methyl-		240	C17H36	006418-43-5	70
3	Decane, 2-methyl-		156	C11H24	006975-98-0	62
4	Nonane, 2-methyl-		142	C10H22	000871-83-0	58
5	Pentadecane, 2,6,10-trimethyl-		254	C18H38	003892-00-0	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122018\
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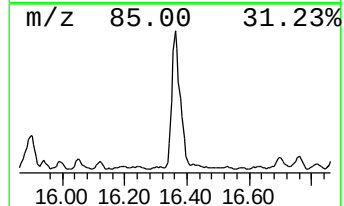
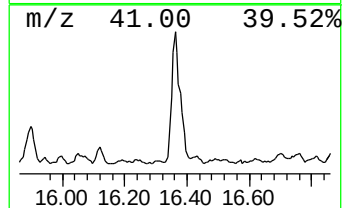
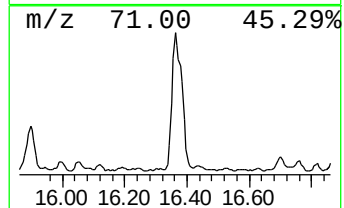
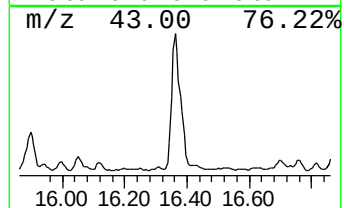
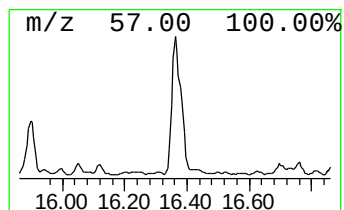
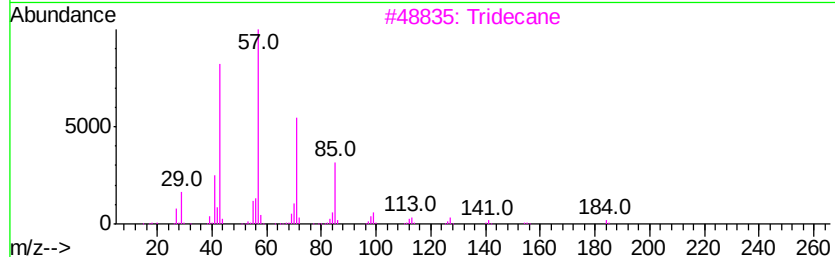
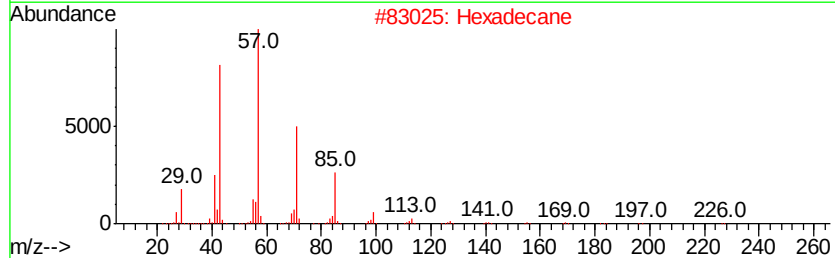
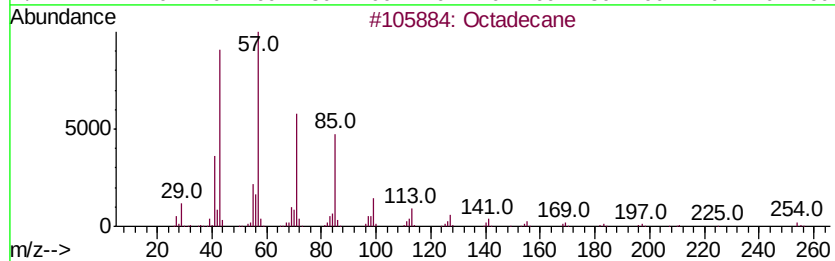
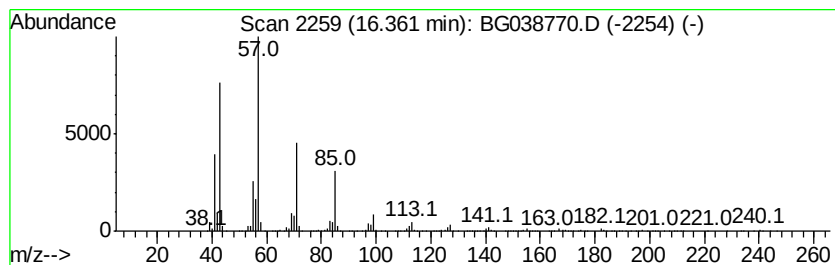
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 Octadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.36	111.80 ng	2300230	Phenanthrene-d10	17.44

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane		254	C18H38	000593-45-3	90
2	Hexadecane		226	C16H34	000544-76-3	90
3	Tridecane		184	C13H28	000629-50-5	86
4	Eicosane, 10-methyl-		296	C21H44	054833-23-7	86
5	Tridecane, 7-hexyl-		268	C19H40	007225-66-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122018\
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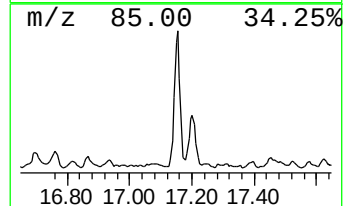
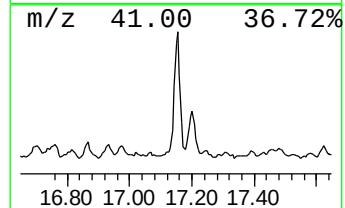
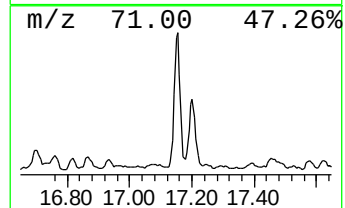
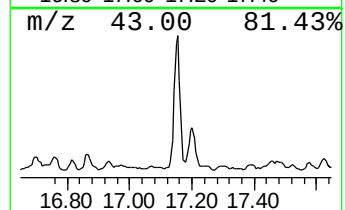
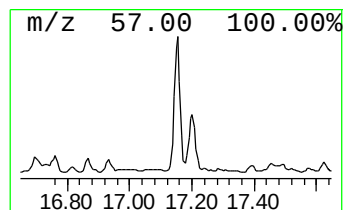
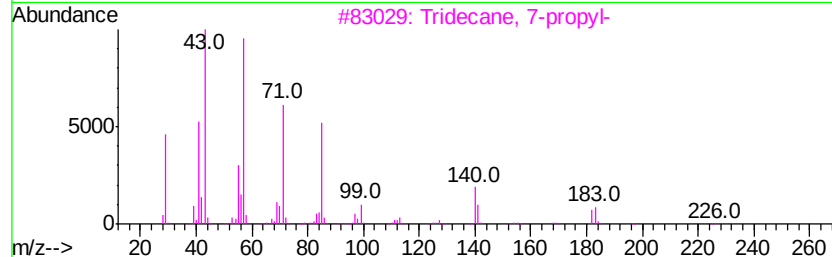
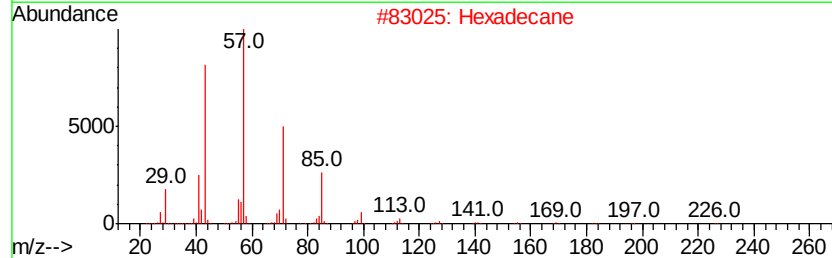
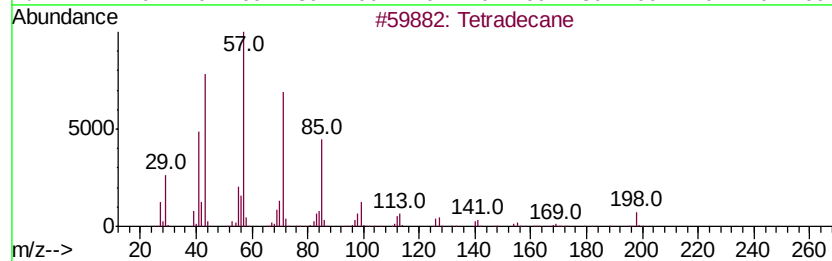
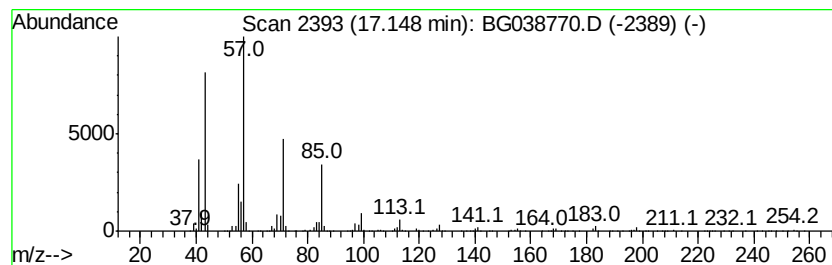
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Tetradecane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.15	43.12 ng	887295	Phenanthrene-d10	17.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	97
2		Hexadecane	226	C16H34	000544-76-3	96
3		Tridecane, 7-propyl-	226	C16H34	055045-09-5	90
4		Tridecane, 6-propyl-	226	C16H34	055045-10-8	87
5		Tridecane	184	C13H28	000629-50-5	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122018\
Data File : BG038770.D
Acq On : 20 Dec 2018 20:12
Operator : JU/SJ
Sample : J6426-08
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P001-SS013-3036-01

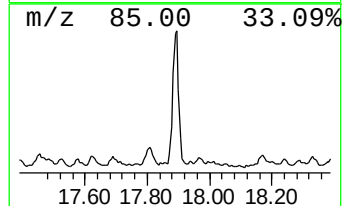
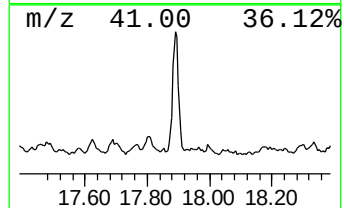
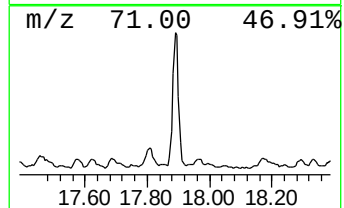
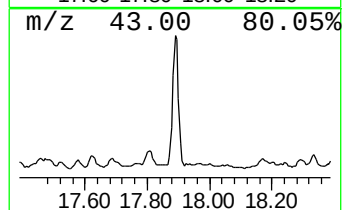
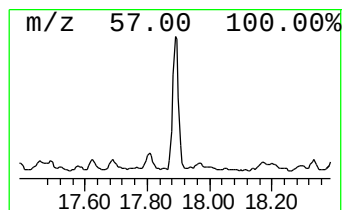
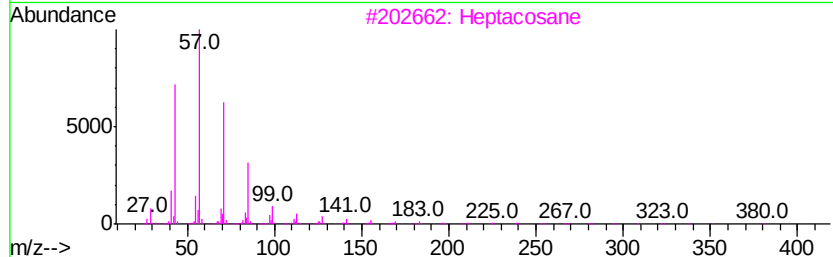
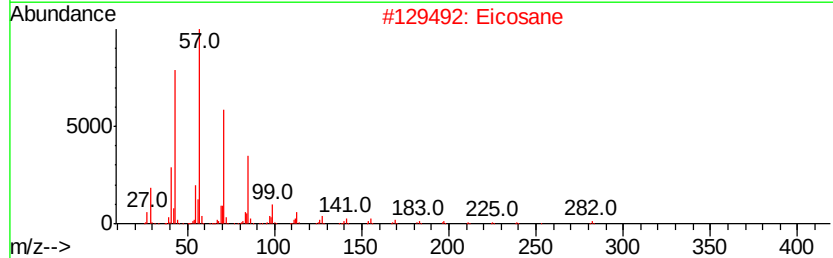
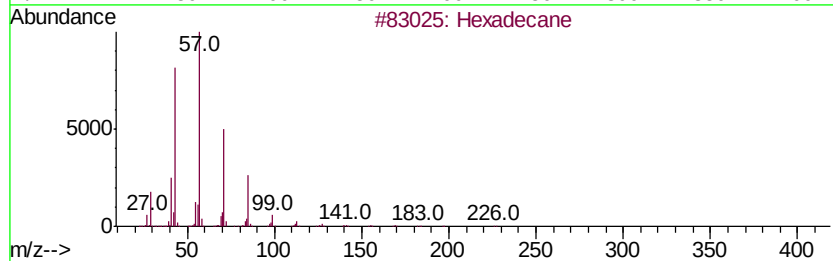
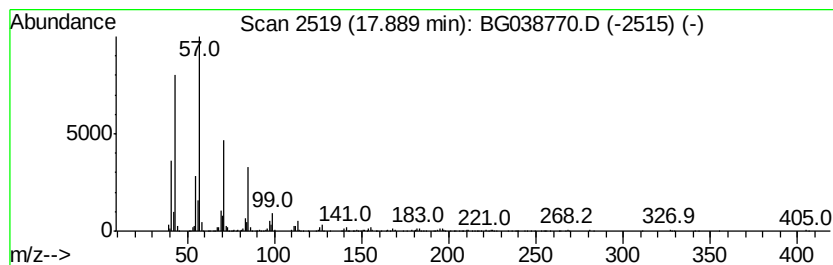
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Hexadecane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.89	41.69 ng	857778	Phenanthrene-d10	17.44

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	94
2			Eicosane	282	C20H42	000112-95-8	91
3			Heptacosane	380	C27H56	000593-49-7	91
4			Tetracosane	338	C24H50	000646-31-1	90
5			Tridecane, 7-propyl-	226	C16H34	055045-09-5	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG122018\
Data File : BG038770.D
Acq On : 20 Dec 2018 20:12
Operator : JU/SJ
Sample : J6426-08
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
P001-SS013-3036-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG121218.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
Benzene, 1,3-dime...	5.70	1264.0	ng	22062600	1	8.06	349102	20.0
unknown7.59	7.59	41.7	ng	727542	1	8.06	349102	20.0
Plumbane, diethyl...	9.56	135.2	ng	2764060	2	10.87	408789	20.0
Dodecane	10.80	61.6	ng	1259510	2	10.87	408789	20.0
Tridecane	12.25	76.2	ng	1557540	2	10.87	408789	20.0
Naphthalene, 1,7-...	13.87	51.6	ng	720403	3	14.69	278998	20.0
Naphthalene, 2,6-...	14.02	77.0	ng	1074720	3	14.69	278998	20.0
Pentadecane	14.55	103.5	ng	1443300	3	14.69	278998	20.0
Z-2-Tridecen-1-ol	15.90	51.5	ng	717913	3	14.69	278998	20.0
Octadecane	16.36	111.8	ng	2300230	4	17.44	411507	20.0
Tetradecane	17.15	43.1	ng	887295	4	17.44	411507	20.0
Hexadecane	17.89	41.7	ng	857778	4	17.44	411507	20.0