

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102918\
 Data File : BG037632.D
 Acq On : 29 Oct 2018 21:06
 Operator : JU/SJ
 Sample : J5665-05
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DL-08A

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101818.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	5.180	317	322	332	rVB	57508	94997	2.88%	0.451%
2	5.639	393	400	410	rBV	856384	1447510	43.85%	6.866%
3	7.248	666	674	684	rBV	934782	1819535	55.12%	8.631%
4	7.624	730	738	750	rBV	875700	1565592	47.43%	7.426%
5	8.100	812	819	833	rVB	198798	359284	10.88%	1.704%
6	8.424	866	874	886	rBV	617954	1114975	33.78%	5.289%
7	9.275	1012	1019	1029	rBV	550314	1031960	31.26%	4.895%
8	10.921	1290	1299	1306	rBV	295591	550434	16.68%	2.611%
9	13.359	1704	1714	1729	rBV	1559431	2564262	77.69%	12.163%
10	14.176	1845	1853	1866	rVB	202749	317378	9.62%	1.505%
11	14.734	1939	1948	1957	rBV2	481922	832967	25.24%	3.951%
12	16.220	2194	2201	2215	rVB	1122502	1763618	53.43%	8.365%
13	17.483	2409	2416	2427	rBV2	626479	1022889	30.99%	4.852%
14	18.335	2556	2561	2569	rBV	130688	221763	6.72%	1.052%
15	19.522	2760	2763	2770	rVB2	25547	40745	1.23%	0.193%
16	19.605	2771	2777	2783	rBV4	47672	74550	2.26%	0.354%
17	19.887	2822	2825	2832	rVB	26039	40214	1.22%	0.191%
18	20.080	2852	2858	2871	rBV	2458401	3300771	100.00%	15.656%
19	20.850	2985	2989	2995	rBV4	20417	33847	1.03%	0.161%
20	21.379	3073	3079	3093	rBV3	114544	318488	9.65%	1.511%
21	21.767	3137	3145	3152	rBV	597758	1087429	32.94%	5.158%
22	21.919	3168	3171	3177	rVB2	29251	39607	1.20%	0.188%
23	22.542	3273	3277	3282	rVB2	31896	54508	1.65%	0.259%
24	23.259	3393	3399	3406	rVB2	46117	99628	3.02%	0.473%
25	24.082	3534	3539	3546	rVB2	52280	119661	3.63%	0.568%
26	25.098	3699	3712	3728	rBV2	304713	1071921	32.47%	5.084%
27	26.226	3898	3904	3913	rVB5	33092	93992	2.85%	0.446%

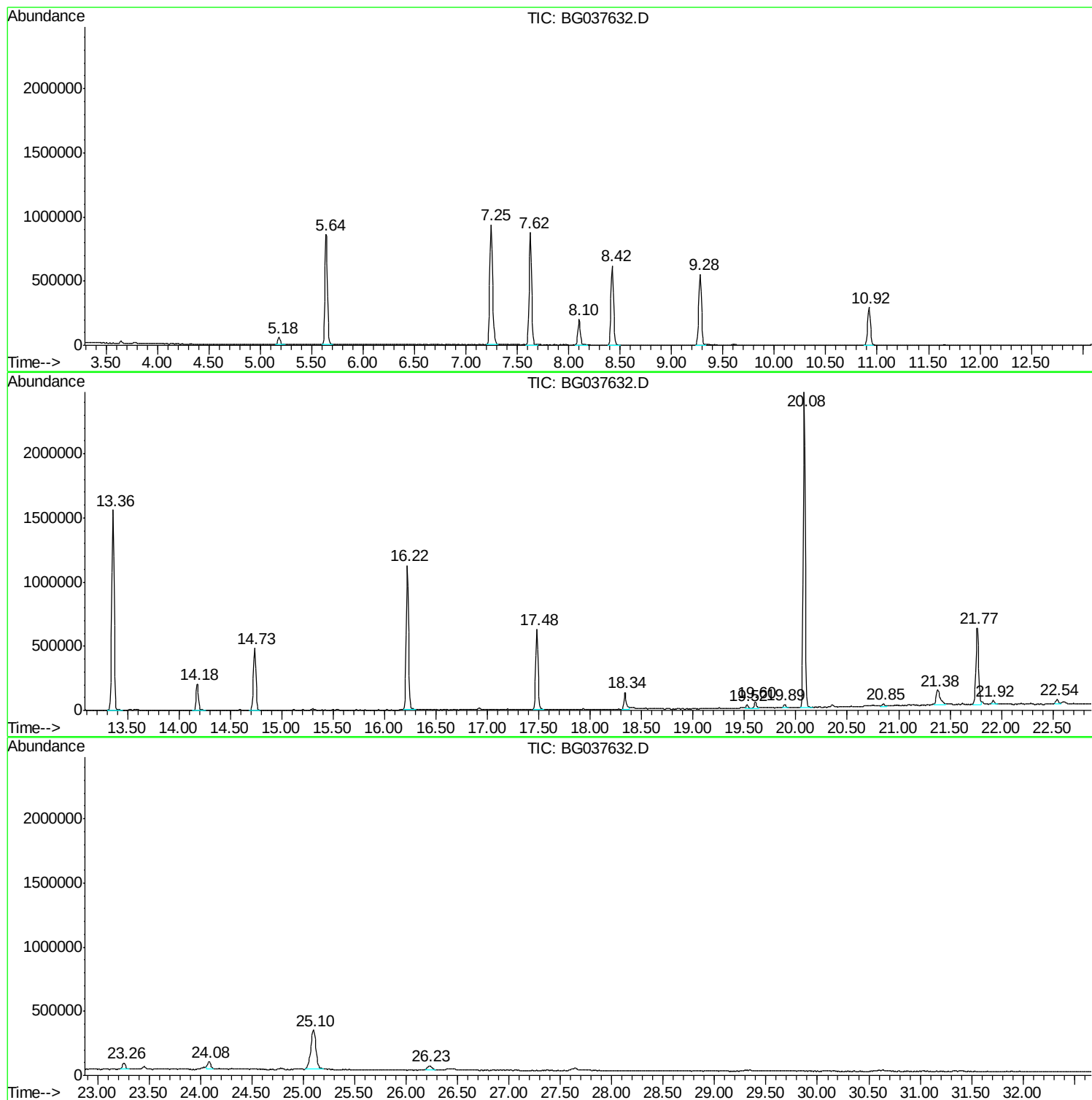
Sum of corrected areas: 21082525

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102918\
Data File : BG037632.D
Acq On : 29 Oct 2018 21:06
Operator : JU/SJ
Sample : J5665-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DL-08A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102918\
Data File : BG037632.D
Acq On : 29 Oct 2018 21:06
Operator : JU/SJ
Sample : J5665-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

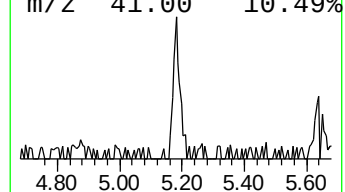
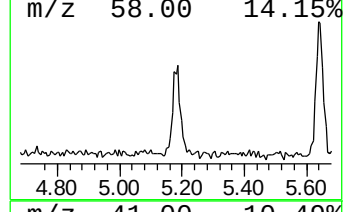
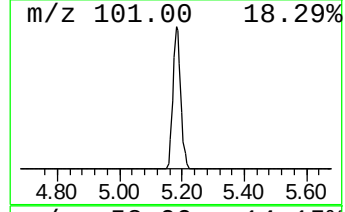
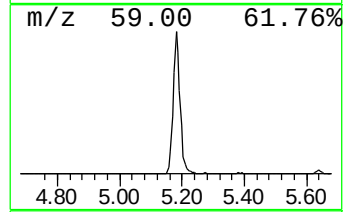
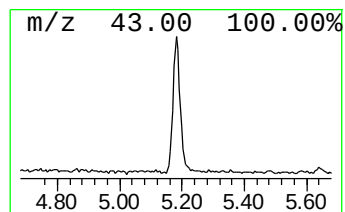
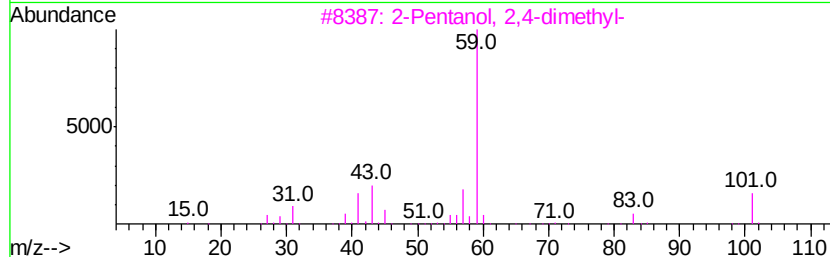
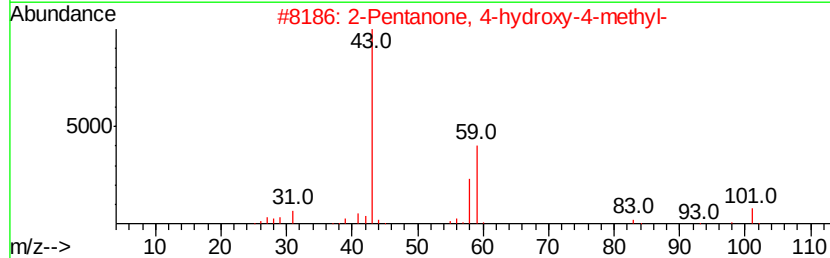
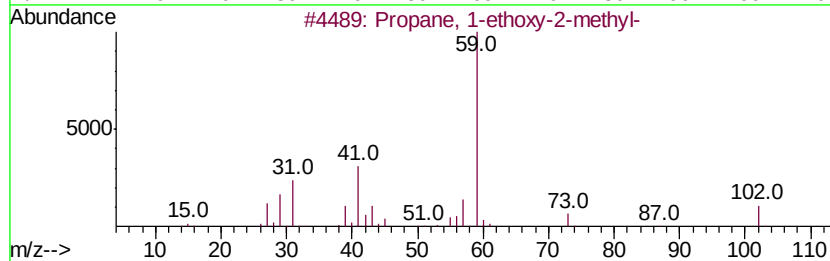
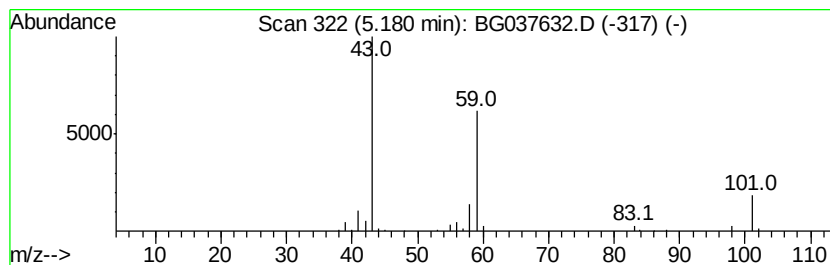
Instrument :
BNA_G
ClientSampleId :
DL-08A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.18	5.29 ng	94997	1,4-Dichlorobenzene-d4	8.10	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	43
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
3	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	38
4	1,2-Butanediol	90	C4H10O2	000584-03-2	28
5	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	23



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102918\
Data File : BG037632.D
Acq On : 29 Oct 2018 21:06
Operator : JU/SJ
Sample : J5665-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DL-08A

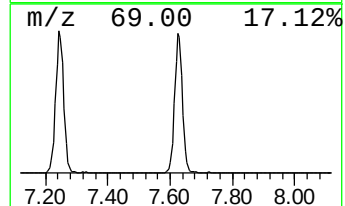
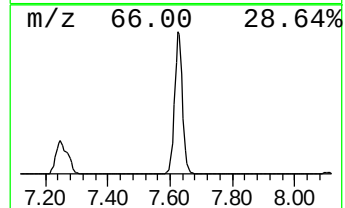
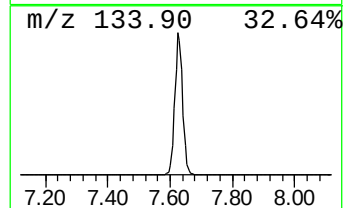
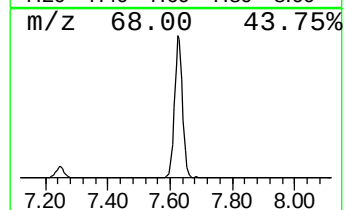
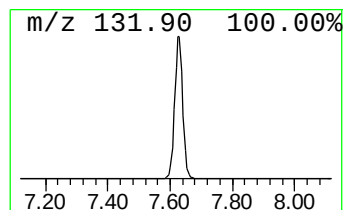
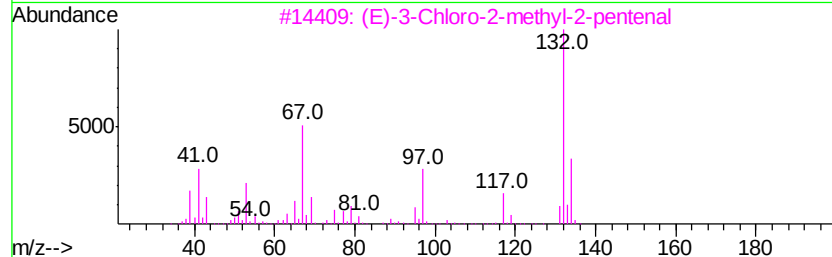
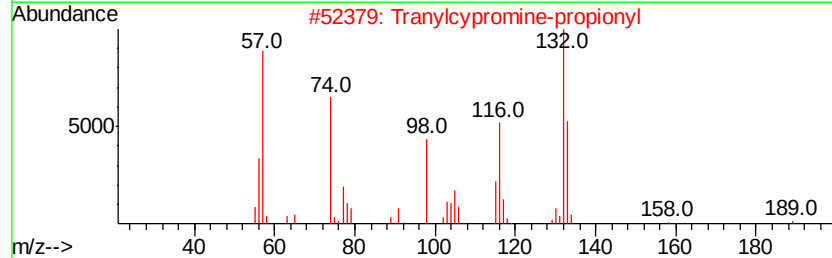
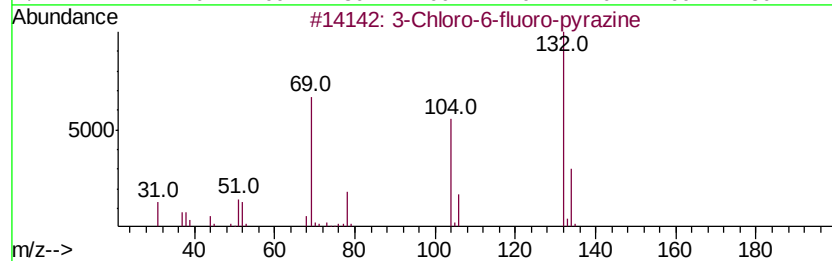
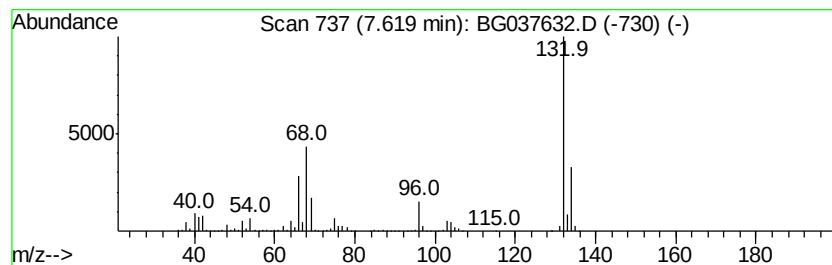
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	87.15 ng	1565590	1,4-Dichlorobenzene-d4	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102918\
Data File : BG037632.D
Acq On : 29 Oct 2018 21:06
Operator : JU/SJ
Sample : J5665-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DL-08A

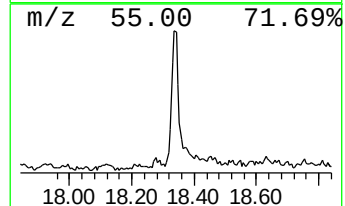
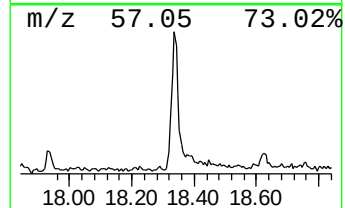
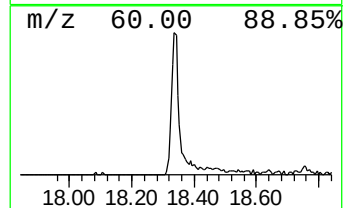
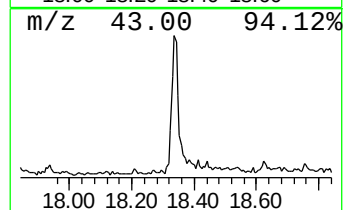
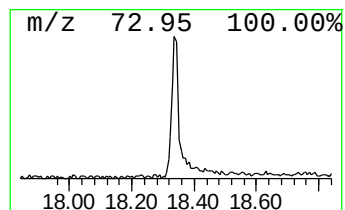
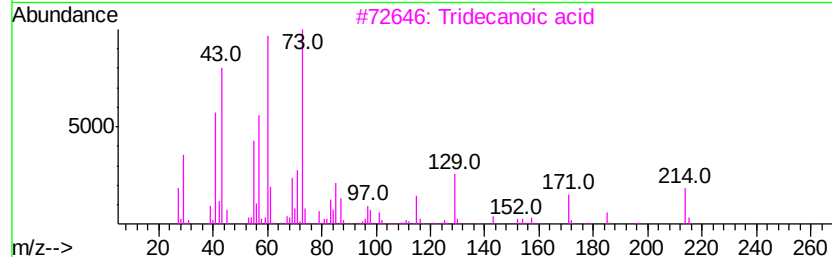
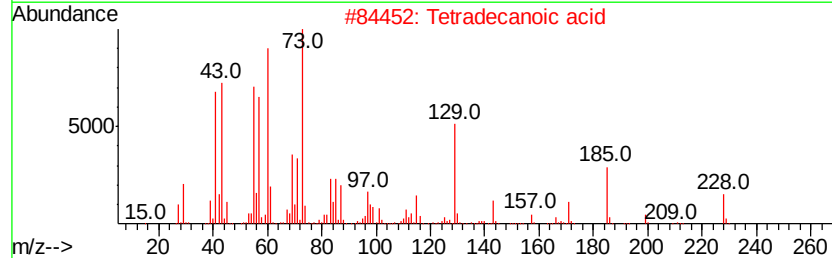
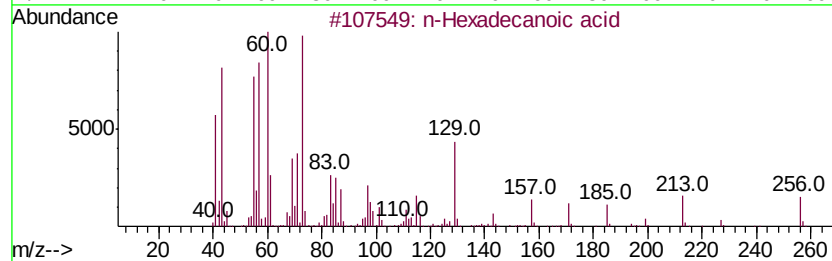
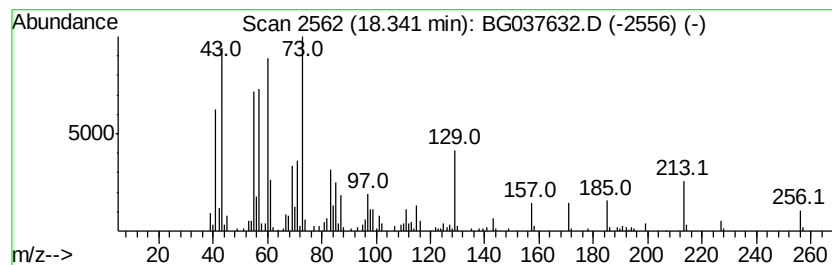
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.34	4.34 ng	221763	Phenanthrene-d10	17.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3	Tridecanoic acid	214	C13H26O2	000638-53-9	81
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	81
5	Dodecanoic acid	200	C12H24O2	000143-07-7	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102918\
Data File : BG037632.D
Acq On : 29 Oct 2018 21:06
Operator : JU/SJ
Sample : J5665-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

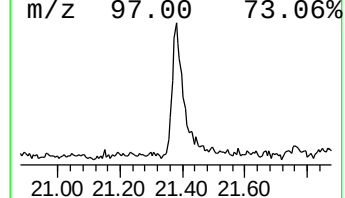
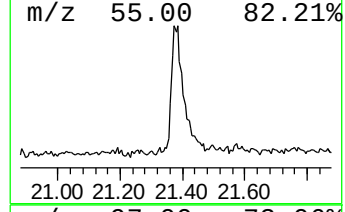
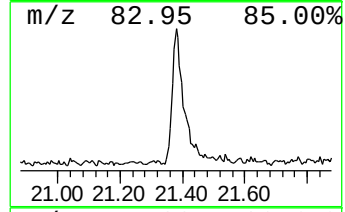
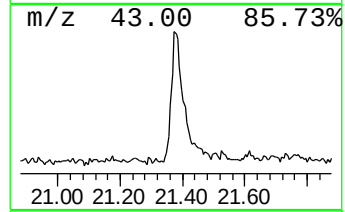
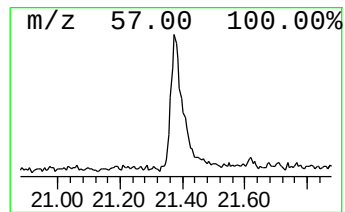
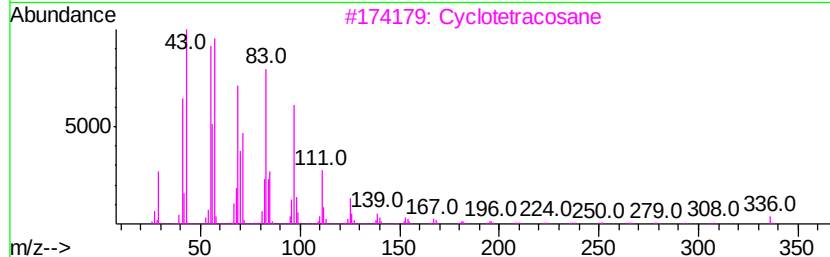
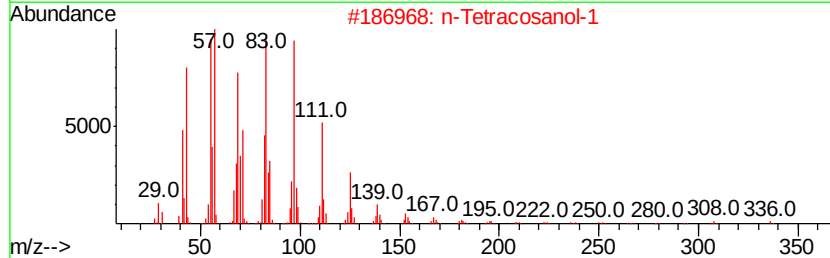
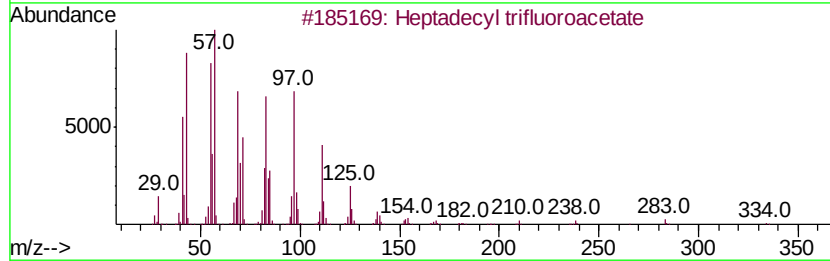
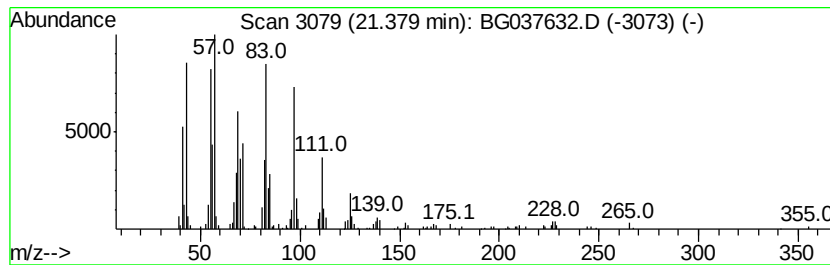
Instrument :
BNA_G
ClientSampleId :
DL-08A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Heptadecyl trifluoroacetate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.		
21.38	5.86 ng	318488	Chrysene-d12		21.77		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	94
2			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
3			Cyclotetracosane	336	C24H48	000297-03-0	91
4			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
5			1-Tricosene	322	C23H46	018835-32-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102918\
Data File : BG037632.D
Acq On : 29 Oct 2018 21:06
Operator : JU/SJ
Sample : J5665-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DL-08A

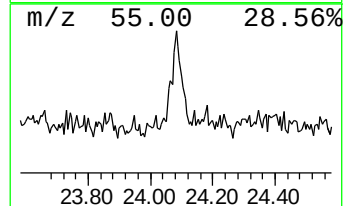
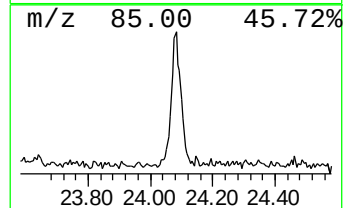
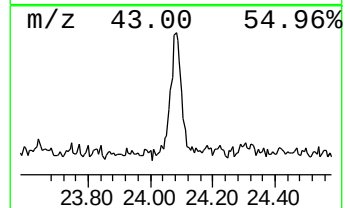
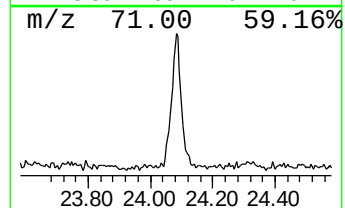
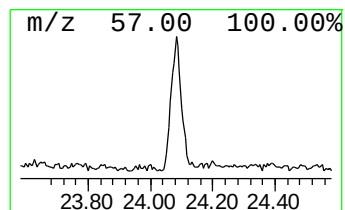
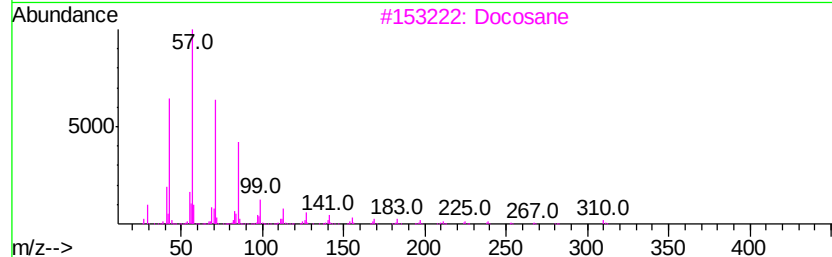
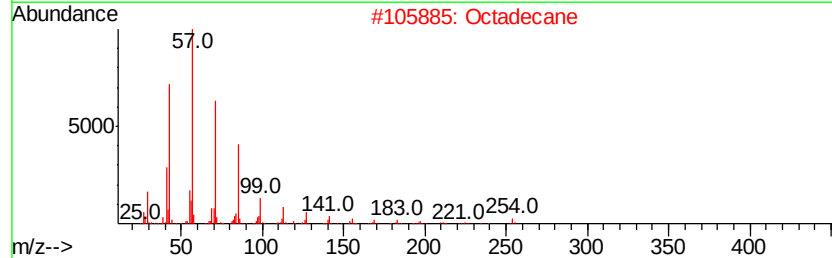
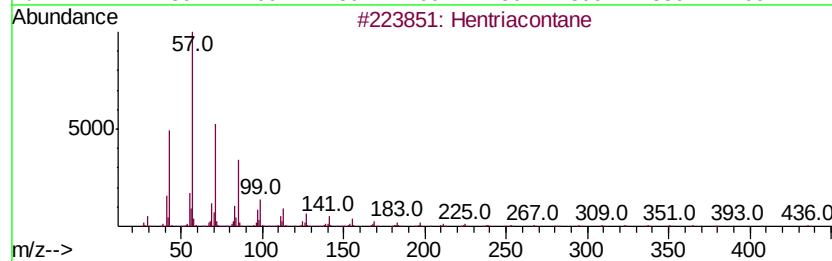
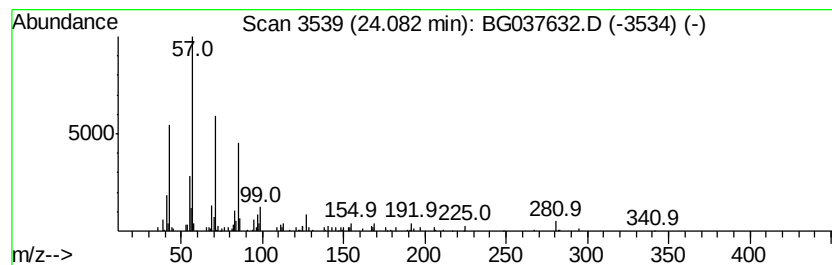
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Hentriacontane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.08	2.23 ng	119661	Perylene-d12	25.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hentriacontane	437	C31H64	000630-04-6	83
2		Octadecane	254	C18H38	000593-45-3	80
3		Docosane	310	C22H46	000629-97-0	80
4		Triacontane	422	C30H62	000638-68-6	80
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG102918\
Data File : BG037632.D
Acq On : 29 Oct 2018 21:06
Operator : JU/SJ
Sample : J5665-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DL-08A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	5.3	ng	94997	1	8.10	359284	20.0
unknown7.62	7.62	87.2	ng	1565590	1	8.10	359284	20.0
n-Hexadecanoic acid	18.34	4.3	ng	221763	4	17.48	1022890	20.0
Heptadecyl triflu...	21.38	5.9	ng	318488	5	21.77	1087430	20.0
Hentriacontane	24.08	2.2	ng	119661	6	25.10	1071920	20.0