

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081420\
 Data File : BG046287.D
 Acq On : 14 Aug 2020 22:43
 Operator : CG/JU
 Sample : L3662-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 VB16182

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-BG073020.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	3.155	7	11	24	rVB	318171	478822	14.00%	1.015%
2	5.535	408	416	426	rBV	1006545	1604435	46.93%	3.401%
3	7.115	677	685	697	rBV	1026727	1754557	51.32%	3.719%
4	7.950	820	827	834	rBV	406306	666077	19.48%	1.412%
5	9.101	1015	1023	1031	rBV	698166	1242078	36.33%	2.633%
6	10.746	1295	1303	1316	rVB	541469	1050827	30.74%	2.228%
7	11.169	1369	1375	1385	rBV6	55307	145402	4.25%	0.308%
8	12.491	1596	1600	1606	rBV5	88713	189414	5.54%	0.402%
9	13.202	1714	1721	1727	rBV	1892516	3059880	89.50%	6.486%
10	13.379	1747	1751	1760	rVB3	352961	633289	18.52%	1.342%
11	13.860	1829	1833	1837	rBV5	278335	435287	12.73%	0.923%
12	14.013	1855	1859	1867	rBV4	519327	1124023	32.88%	2.383%
13	14.424	1926	1929	1936	rVB2	993930	1568314	45.87%	3.324%
14	14.577	1947	1955	1960	rBV4	1153194	3152547	92.21%	6.683%
15	15.382	2088	2092	2097	rBV2	1525598	2483766	72.65%	5.265%
16	16.075	2206	2210	2213	rBV	1393830	2038045	59.61%	4.320%
17	21.569	3140	3145	3151	rBV3	1189686	2948488	86.24%	6.250%
18	21.622	3151	3154	3167	rVB	1155818	2121173	62.04%	4.496%
19	22.380	3278	3283	3288	rBV4	492638	1031256	30.16%	2.186%
20	22.438	3288	3293	3299	rVB2	1855272	2983177	87.25%	6.324%
21	23.702	3501	3508	3515	rBV	804484	2459225	71.93%	5.213%
22	23.825	3524	3529	3534	rBV	912574	1801382	52.69%	3.819%
23	23.878	3534	3538	3545	rVB	523980	1022104	29.90%	2.167%
24	24.395	3621	3626	3639	rVB	492699	1319834	38.60%	2.798%
25	24.536	3643	3650	3660	rVB	550509	1359089	39.75%	2.881%
26	24.683	3668	3675	3681	rBV2	530767	1349262	39.46%	2.860%
27	25.834	3863	3871	3881	rVV2	752747	2188747	64.02%	4.640%
28	25.946	3883	3890	3909	rVB5	441055	1545620	45.21%	3.276%
29	29.830	4539	4551	4575	rVB10	660887	3418973	100.00%	7.247%

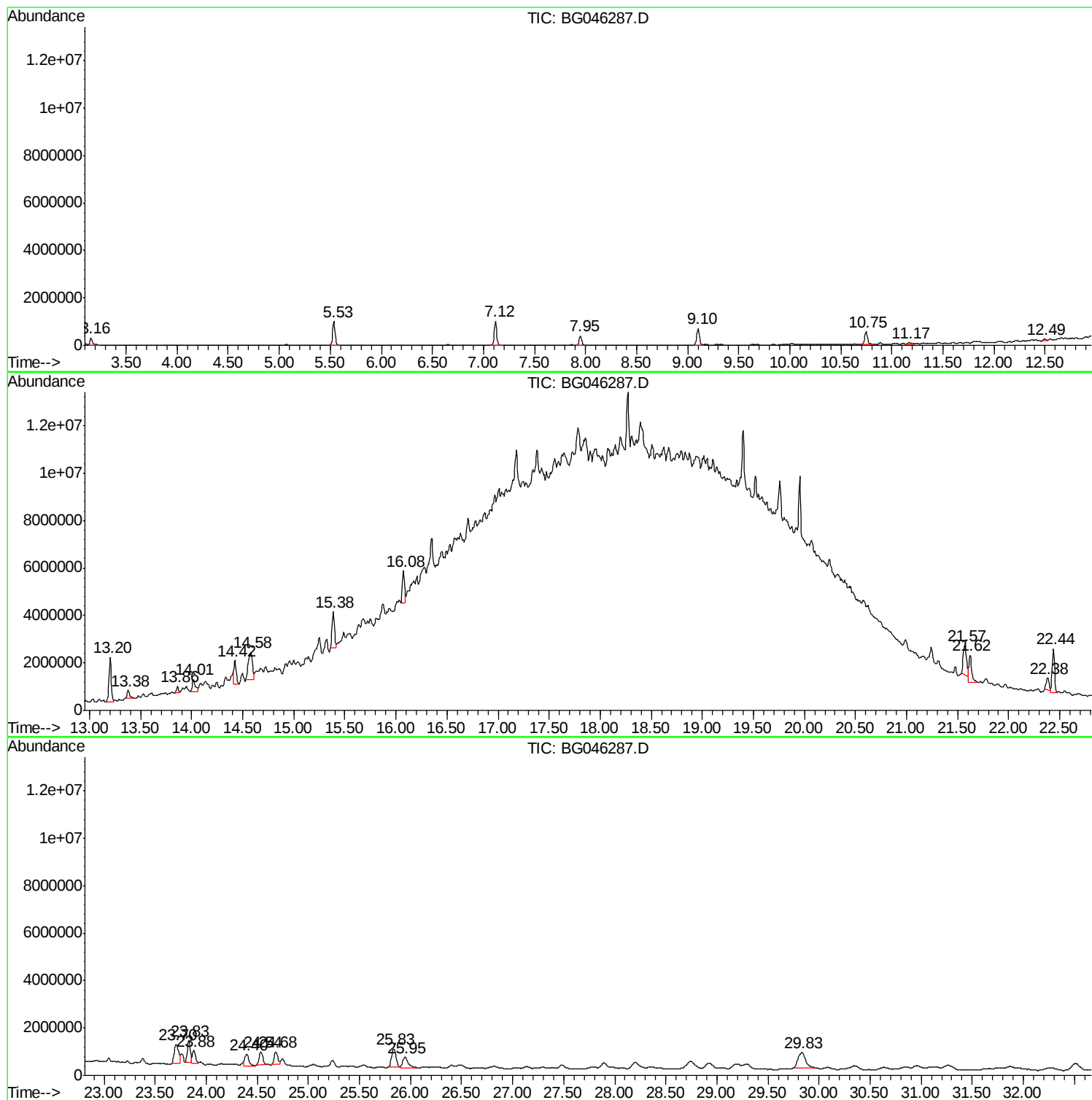
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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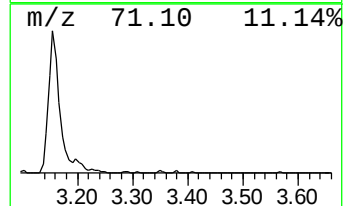
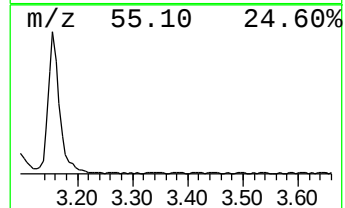
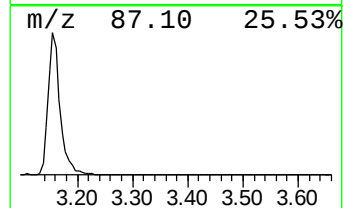
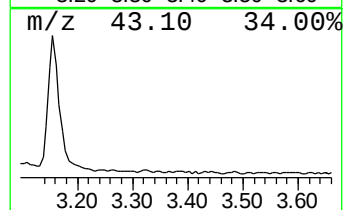
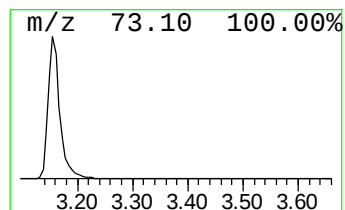
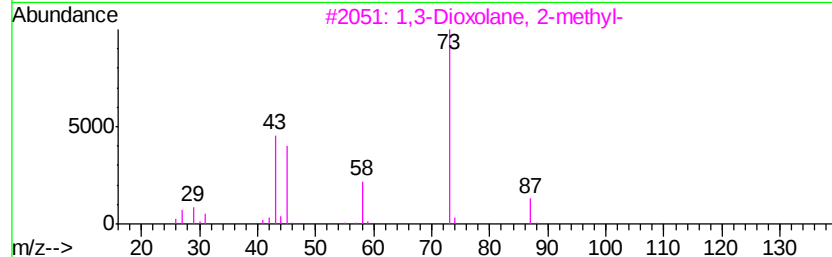
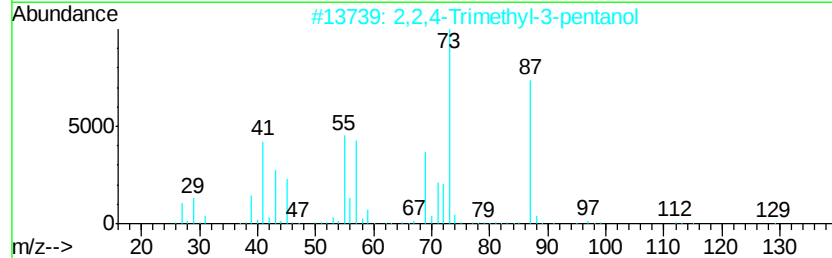
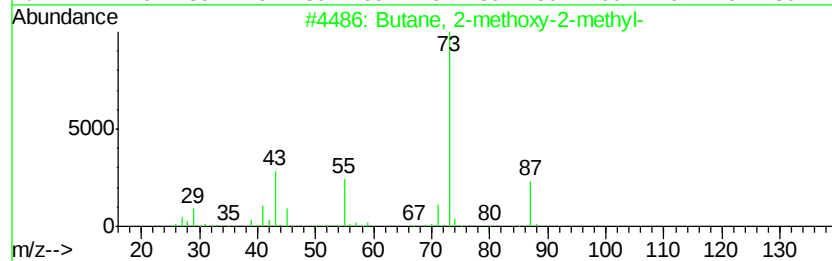
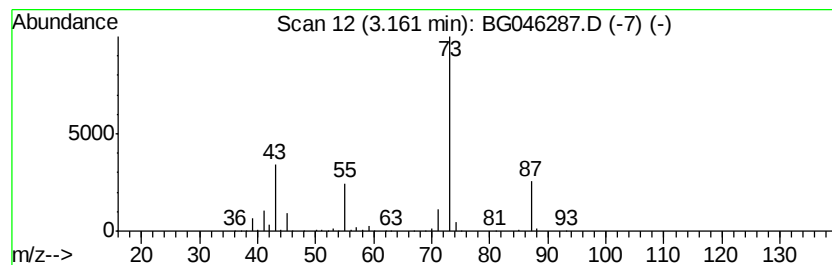
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.16	14.38 ng	478822	1,4-Dichlorobenzene-d4	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	28
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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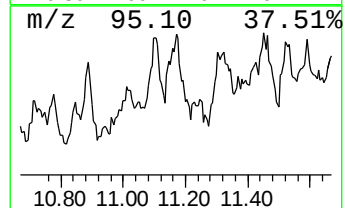
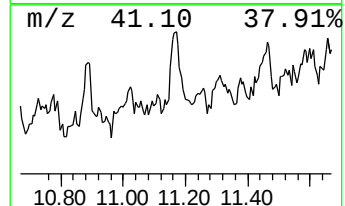
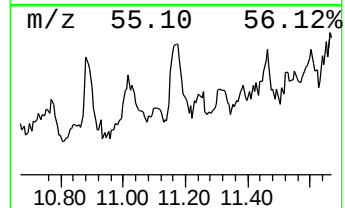
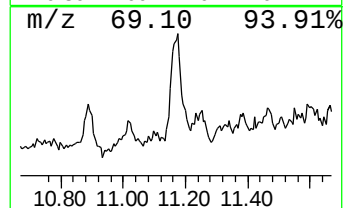
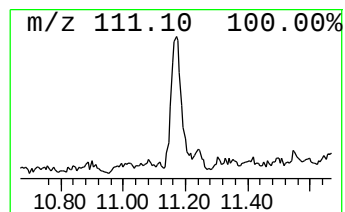
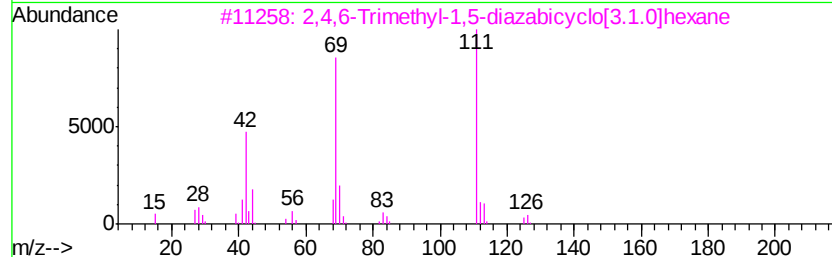
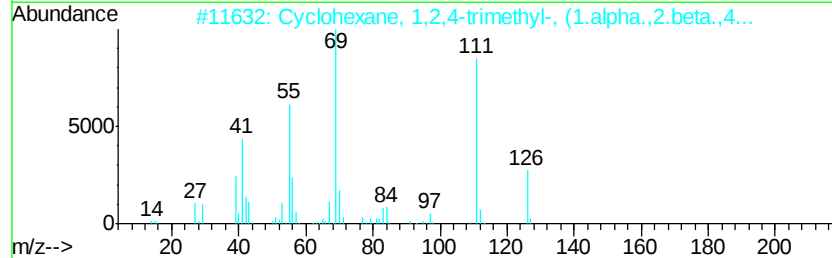
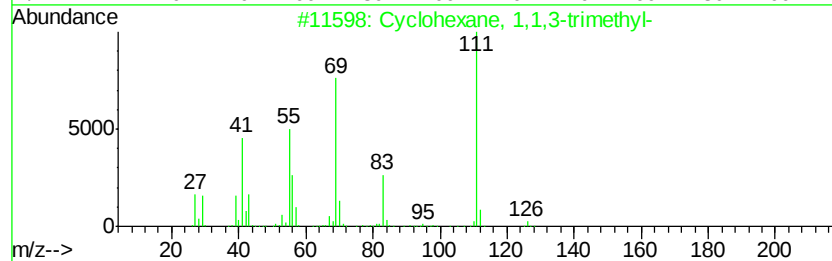
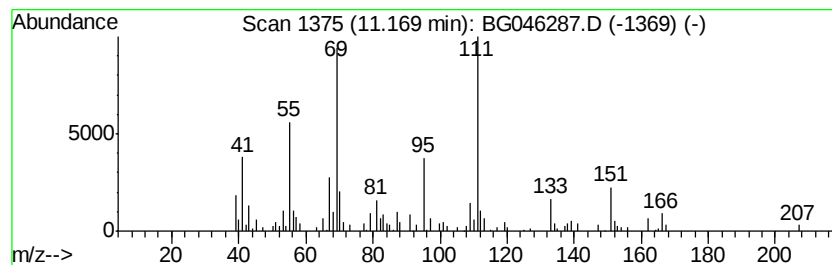
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Peak Number 2 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.17	2.77 ng	145402	Naphthalene-d8	10.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	53
2		Cyclohexane, 1,2,4-trimethyl-, (...)	126	C9H18	007667-60-9	53
3		2,4,6-Trimethyl-1,5-diazabicyclo...	126	C7H14N2	1000283-21-2	53
4		Cyclopentanecarboxamide, N-(4-fl...	207	C12H14FNO	1000307-02-4	50
5		Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	50



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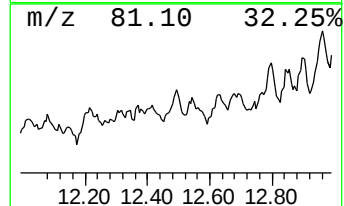
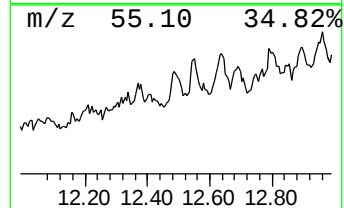
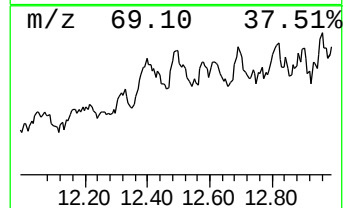
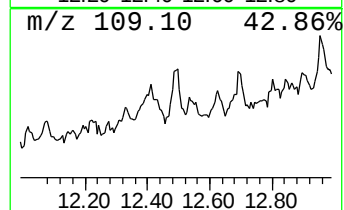
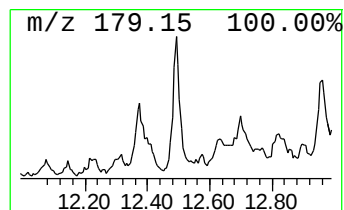
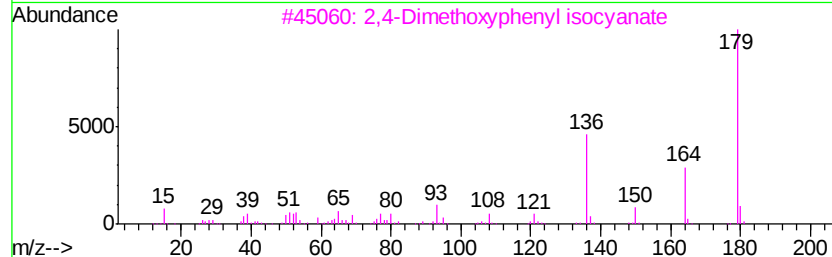
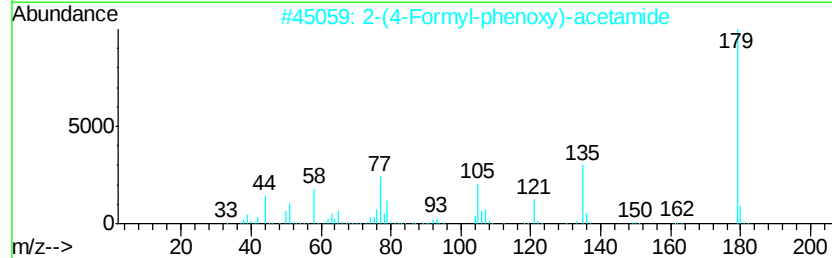
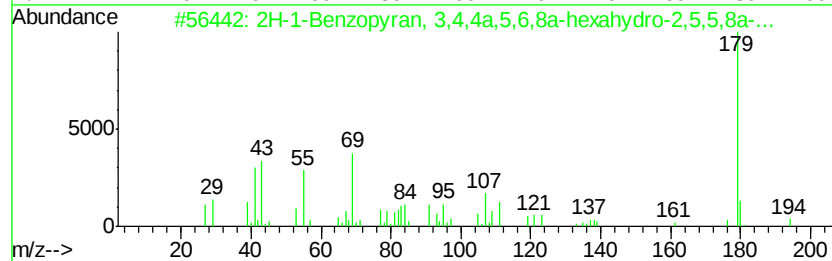
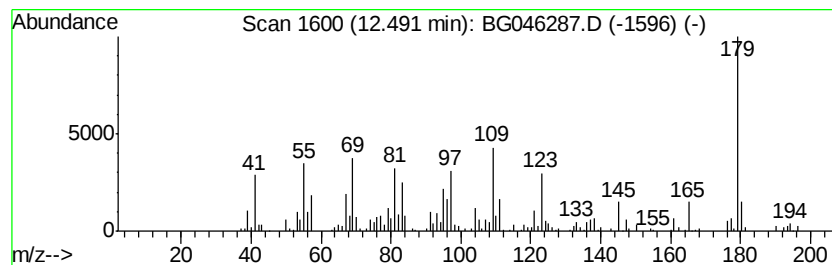
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 unknown12.49 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.49	3.61 ng	189414	Naphthalene-d8	10.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-1-Benzopyran, 3,4,4a,5,6,8a-h...	194	C13H22O	041678-32-4	46
2		2-(4-Formyl-phenoxy)-acetamide	179	C9H9NO3	135857-20-4	38
3		2,4-Dimethoxyphenyl isocyanate	179	C9H9NO3	084370-87-6	38
4		DL-2-fluorophenylglycine, N-dime...	238	C12H15FN2O2	1000375-81-1	37
5		Isoxazole, 3,4-dimethyl-5-(2-thi...	179	C9H9NOS	056421-65-9	35



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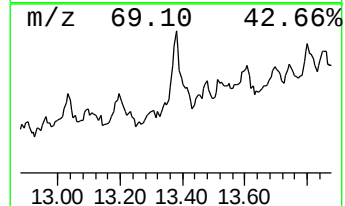
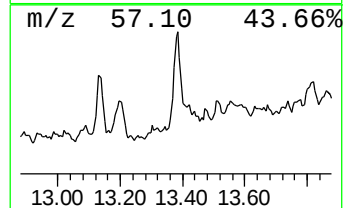
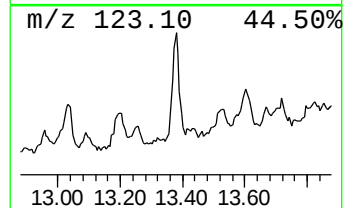
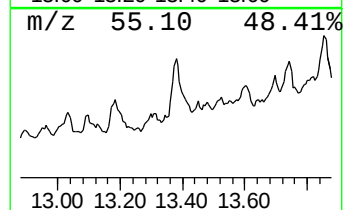
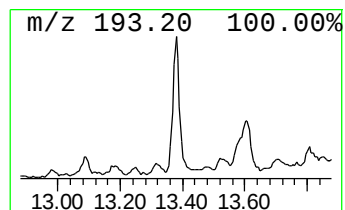
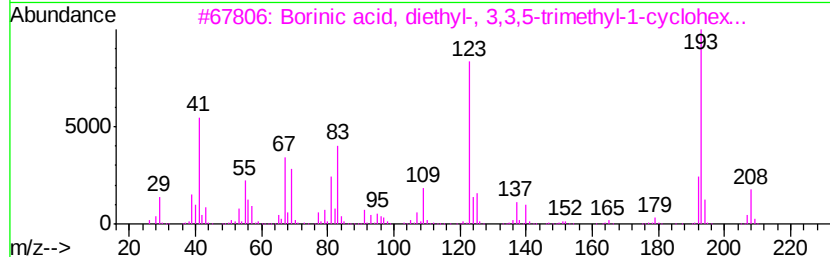
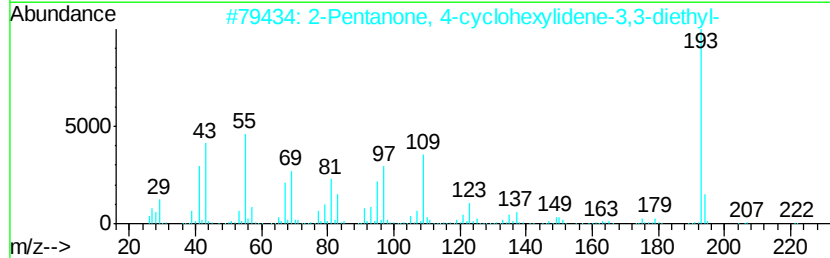
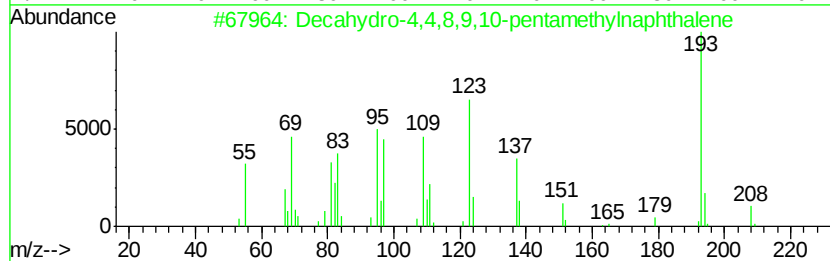
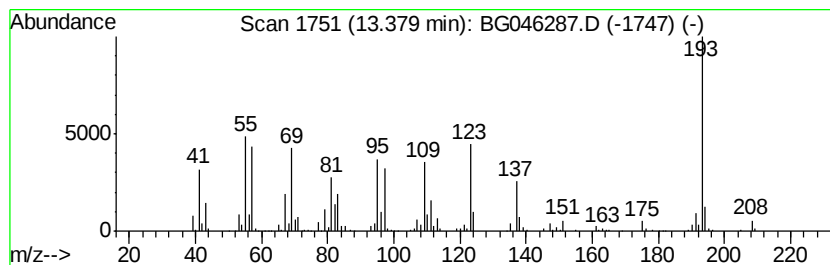
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Peak Number 4 Decahydro-4,4,8,9,10-pentam... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.38	4.02 ng	633289	Acenaphthene-d10	14.58	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	76
2	2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	49
3	Borinic acid, diethyl-, 3,3,5-tr...	208	C13H25BO	057387-76-5	40
4	S-Phenyl 5-(phenylthio)pentaneth...	302	C17H18OS2	1000234-40-7	40
5	Iminostilbene	193	C14H11N	000256-96-2	35



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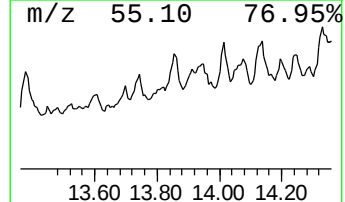
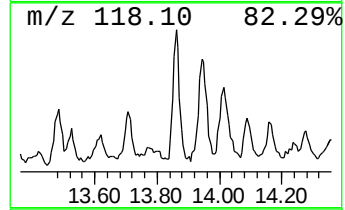
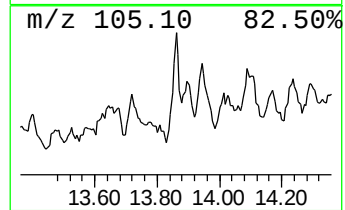
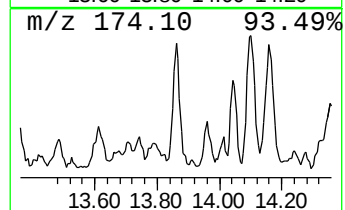
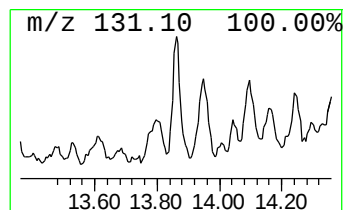
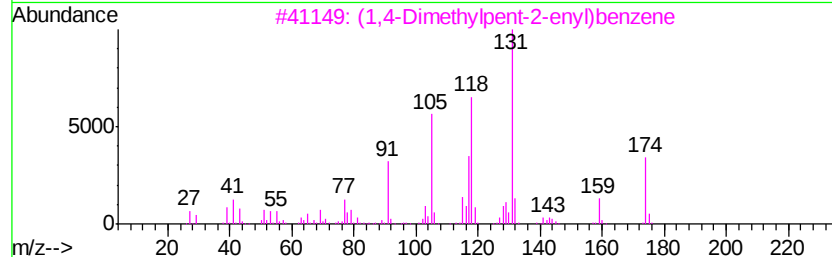
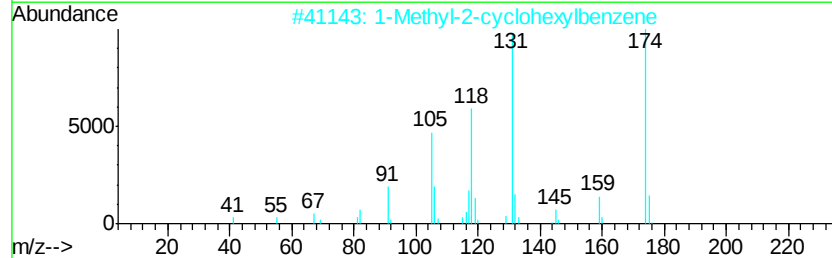
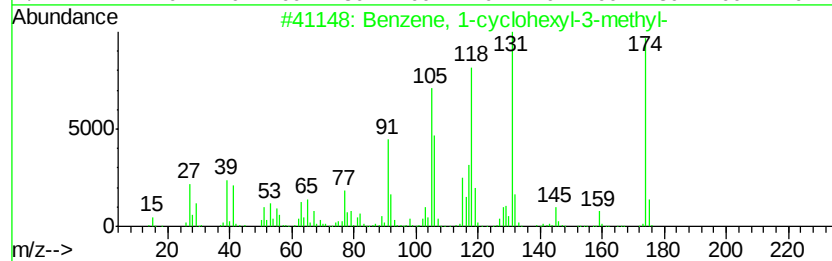
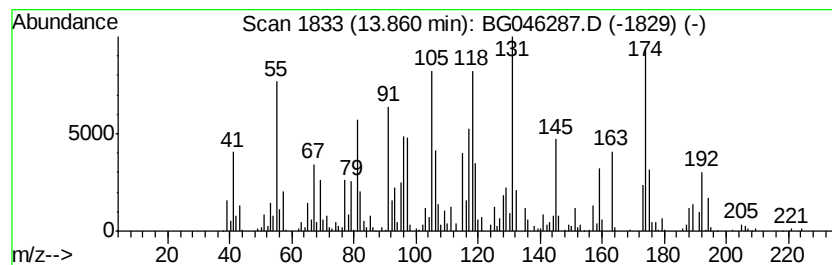
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Peak Number 5 Benzene, 1-cyclohexyl-3-met... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	2.76 ng	435287	Acenaphthene-d10	14.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-cyclohexyl-3-methyl-	174 C13H18	004575-46-6 90
2		1-Methyl-2-cyclohexylbenzene	174 C13H18	004501-35-3 53
3		(1,4-Dimethylpent-2-enyl)benzene	174 C13H18	1000184-97-9 46
4		4-Benzyl-isoxazolidine	163 C10H13NO	1000186-97-4 25
5		2,3-Dihydro-4-methyl-1H-1,5-benz...	174 C10H10N2O	006276-48-8 22



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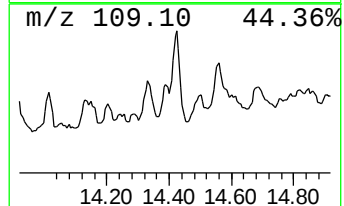
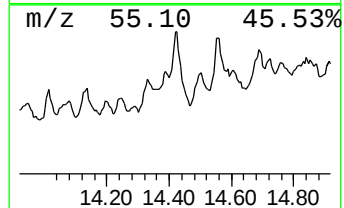
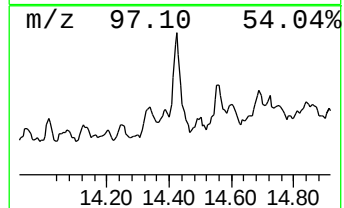
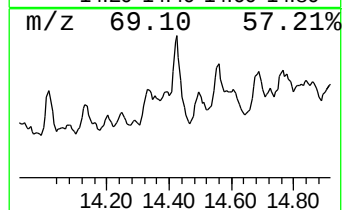
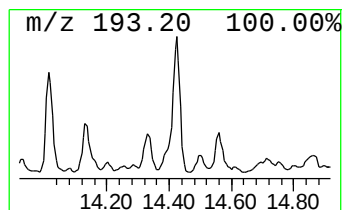
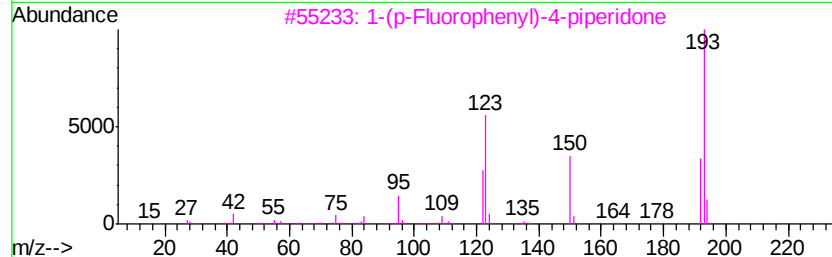
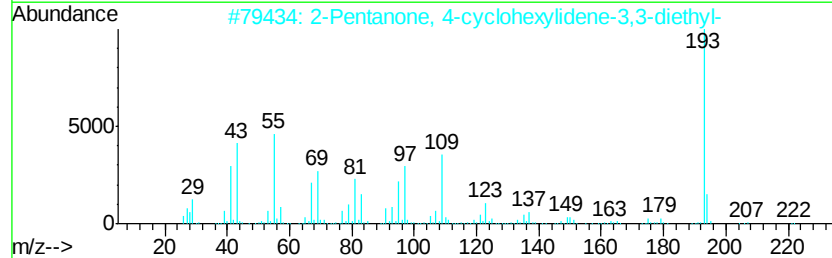
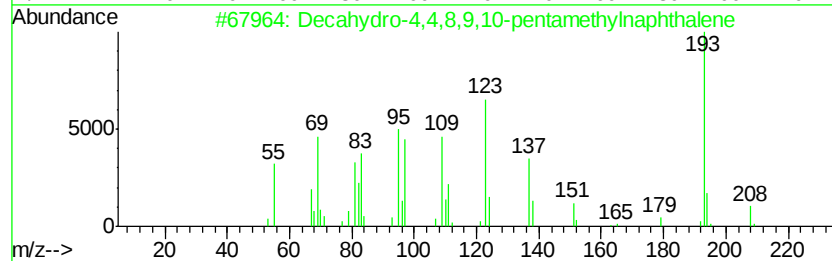
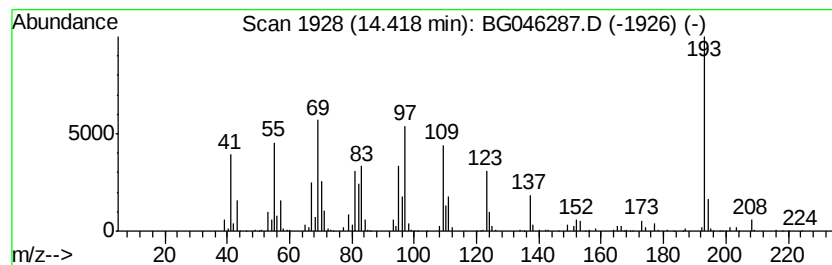
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2-Pentanone, 4-cyclohexylid... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.42	9.95 ng	1568310	Acenaphthene-d10	14.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	53
2		2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	50
3		1-(p-Fluorophenyl)-4-piperidone	193	C11H12FN0	1000238-56-7	32
4		Borinic acid, diethyl-, 3,3,5-tr...	208	C13H25B0	057387-76-5	25
5		2-Amino-4-ethylthiomethyl-6-pipe...	253	C11H19NS	022362-22-7	16



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081420\
Data File : BG046287.D
Acq On : 14 Aug 2020 22:43
Operator : CG/JU
Sample : L3662-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
VB16182

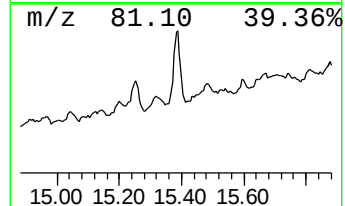
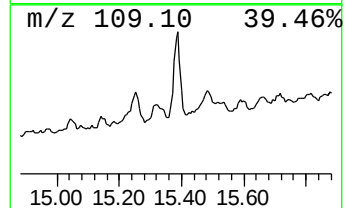
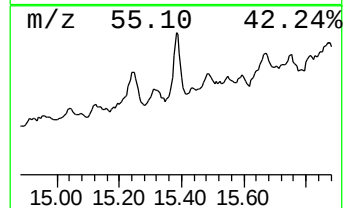
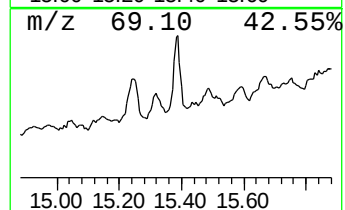
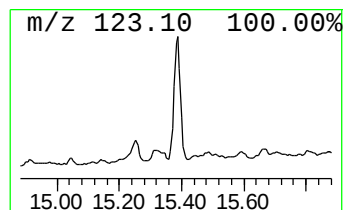
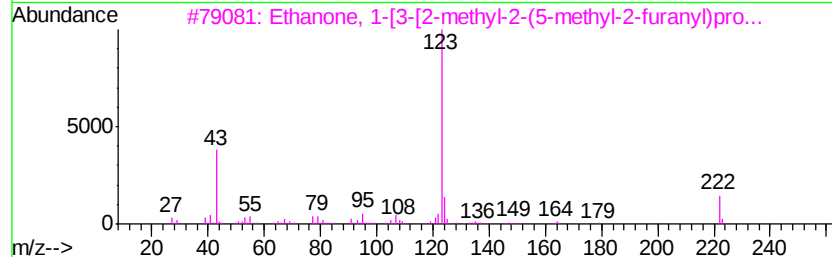
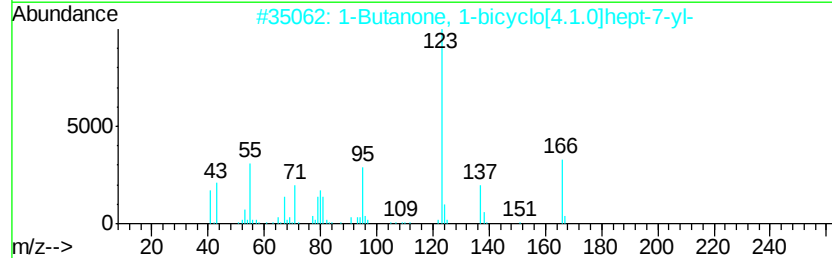
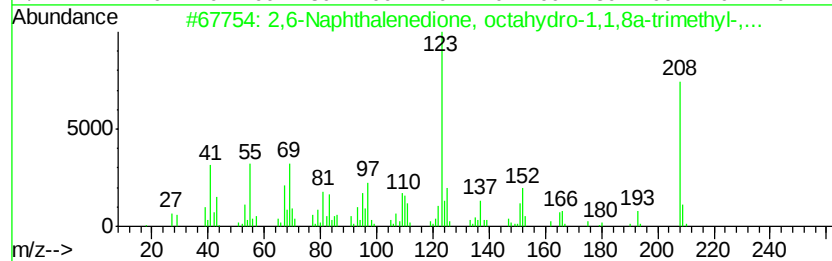
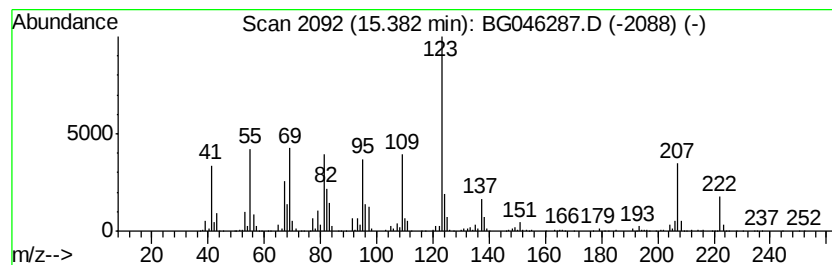
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 2,6-Naphthalenedione, octah... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.38	15.76 ng	2483770	Acenaphthene-d10	14.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,6-Naphthalenedione, octahydro-...	208	C13H20O2	057289-17-5	50
2		1-Butanone, 1-bicyclo[4.1.0]hept...	166	C11H18O	054764-62-4	43
3		Ethanone, 1-[3-[2-methyl-2-(5-me...	222	C13H18O3	080114-28-9	43
4		2,2-Dimethylocta-3,4-dienal	152	C10H16O	000590-71-6	38
5		Imidazolidin-2-one, 1-(2-fluorob...	208	C10H9FN2O2	1000310-62-2	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081420\
 Data File : BG046287.D
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Instrument :
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 ClientSampleId :
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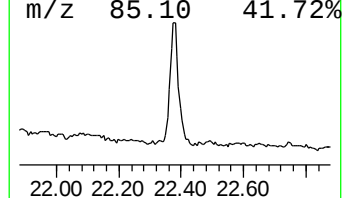
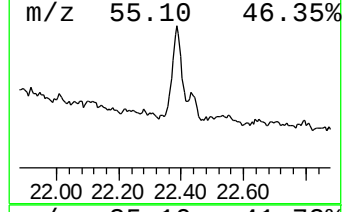
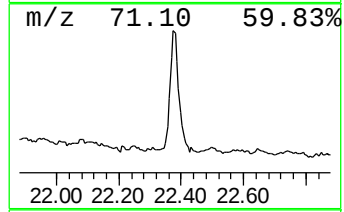
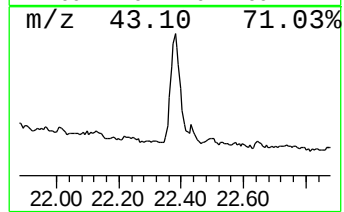
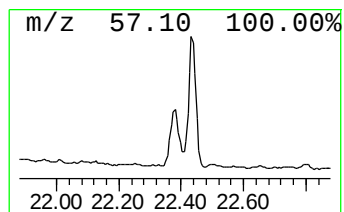
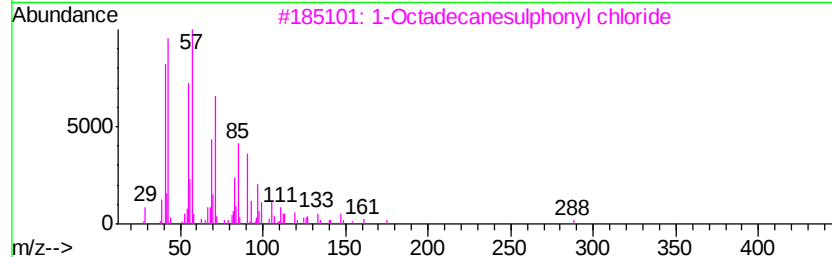
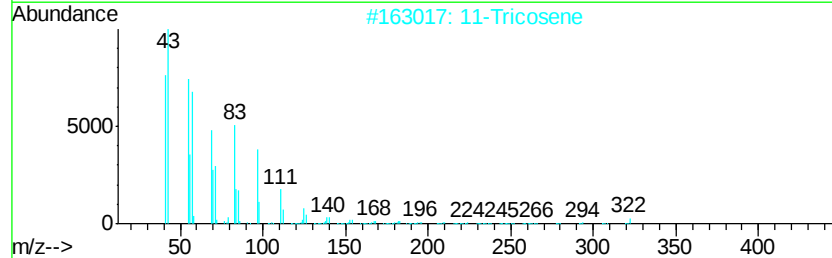
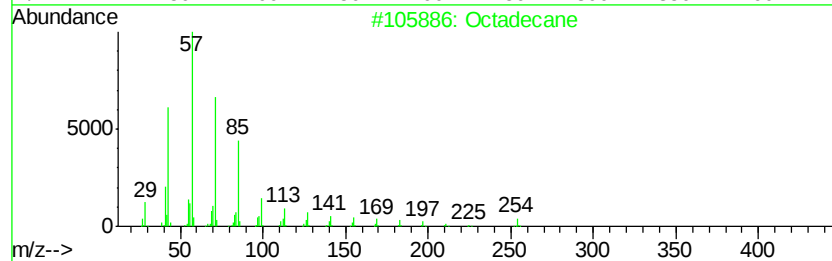
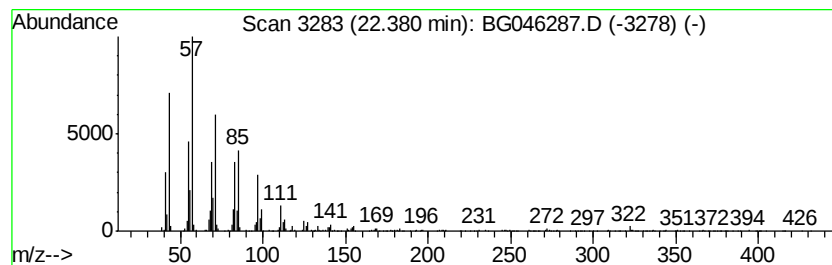
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 8 Octadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.38	7.00 ng	1031260	Chrysene-d12	21.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane		254	C18H38	000593-45-3	93
2	11-Tricosene		322	C23H46	052078-56-5	89
3	1-Octadecanesulphonyl chloride		352	C18H37ClO2S	1000342-70-4	87
4	Octacosane		394	C28H58	000630-02-4	86
5	Tricosane		324	C23H48	000638-67-5	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081420\
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Operator : CG/JU
Sample : L3662-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

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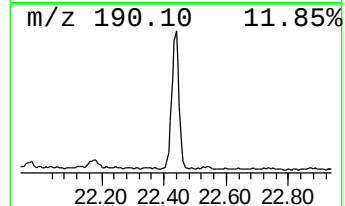
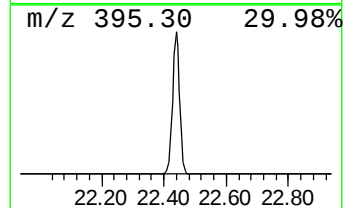
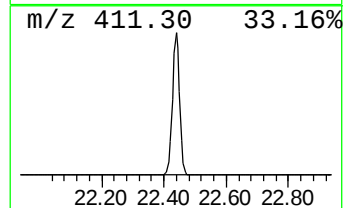
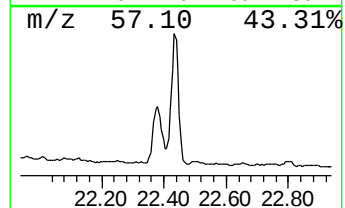
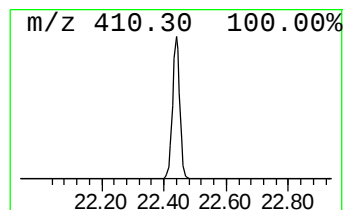
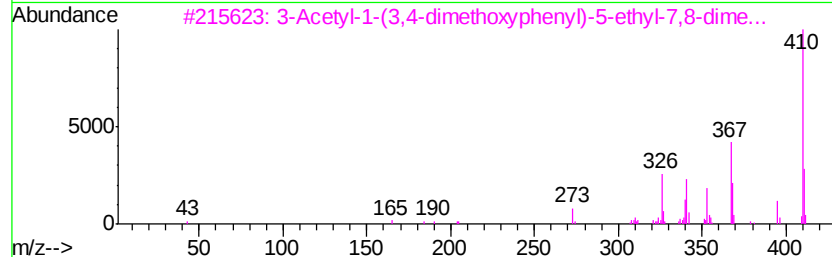
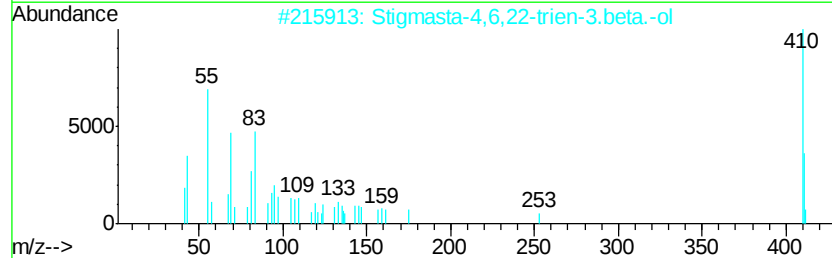
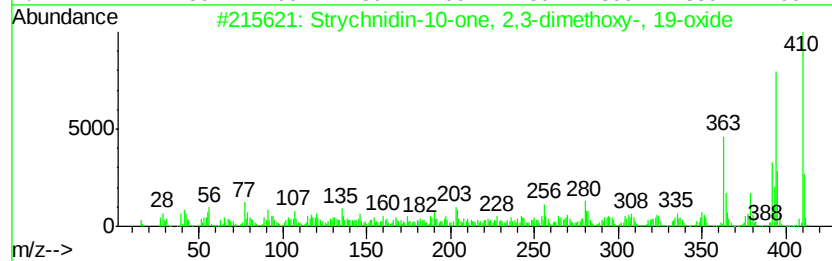
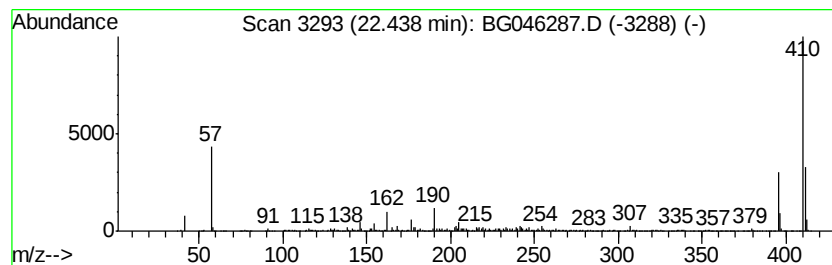
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Strychnidin-10-one, 2,3-dim... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.44	20.24 ng	2983180	Chrysene-d12	21.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Strychnidin-10-one, 2,3-dimethox...	410	C23H26N2O5	017301-81-4	72
2		Stigmasta-4,6,22-trien-3.beta.-ol	410	C29H46O	096737-58-5	58
3		3-Acetyl-1-(3,4-dimethoxyphenyl)...	410	C23H26N2O5	090140-65-1	53
4		Ferrocene, 1,1',3,3'-tetrakis(1,...	410	C26H42Fe	001317-17-5	50
5		24(S)-Ethyl-3.alpha.,5.alpha.-cy...	410	C29H46O	1000105-21-7	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081420\
Data File : BG046287.D
Acq On : 14 Aug 2020 22:43
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Misc :
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ClientSampleId :
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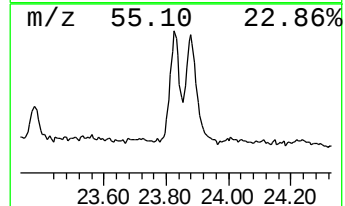
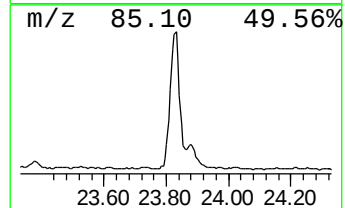
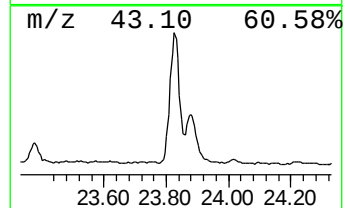
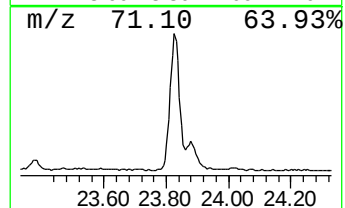
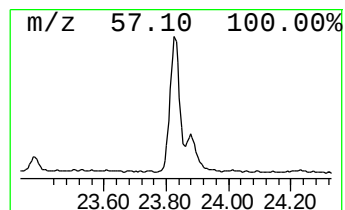
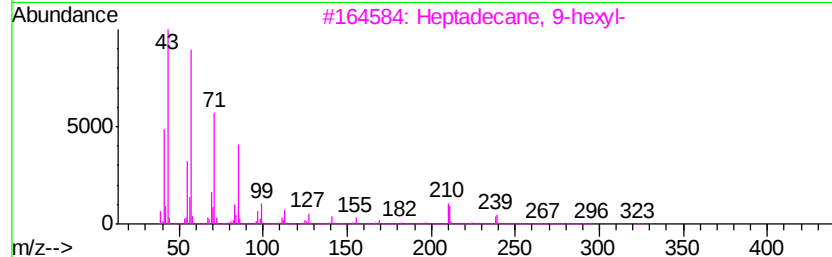
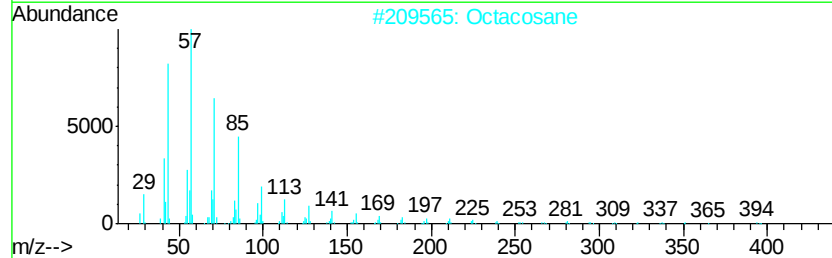
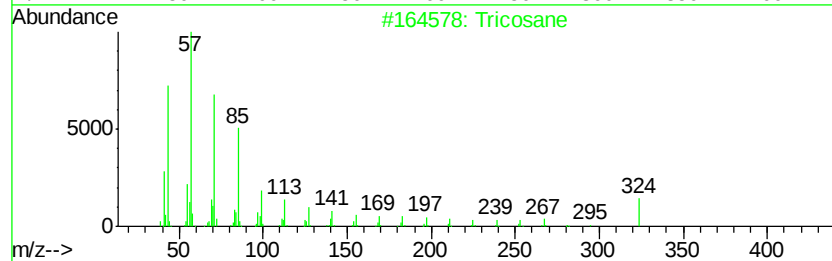
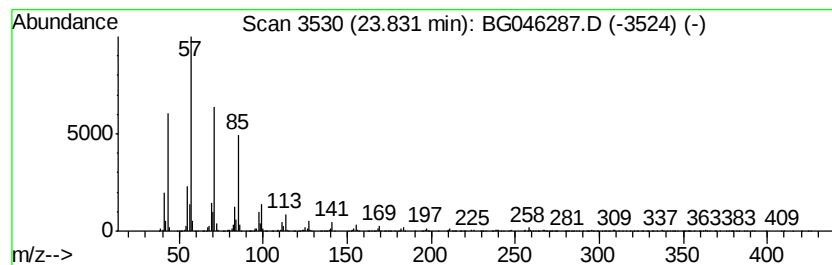
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tricosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.83	26.70 ng	1801380	Perylene-d12	24.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosane	324	C23H48	000638-67-5	97
2		Octacosane	394	C28H58	000630-02-4	97
3		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
4		Hexacosane	366	C26H54	000630-01-3	91
5		Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081420\
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Misc :
ALS Vial : 14 Sample Multiplier: 1

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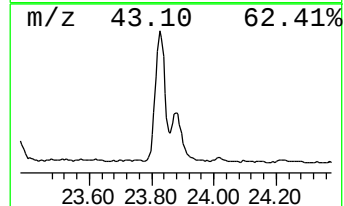
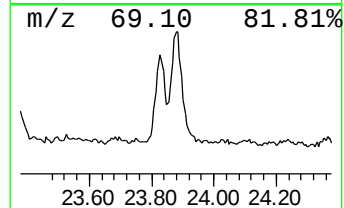
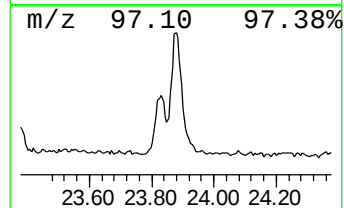
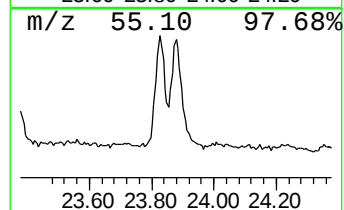
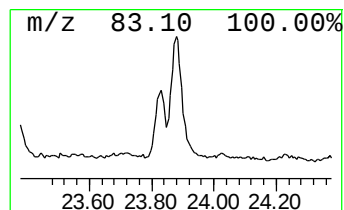
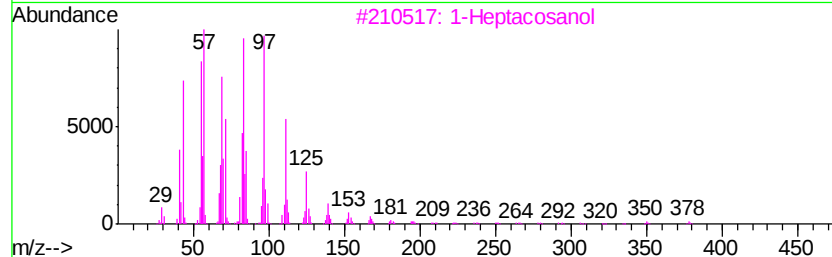
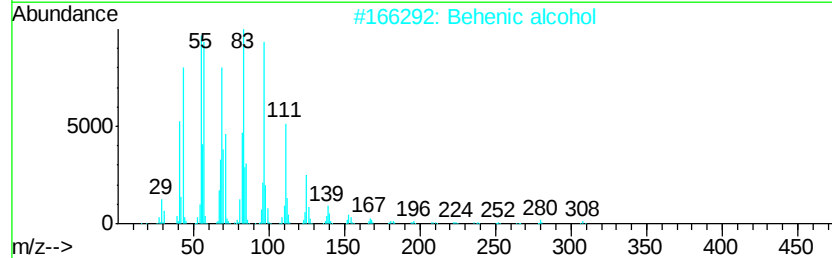
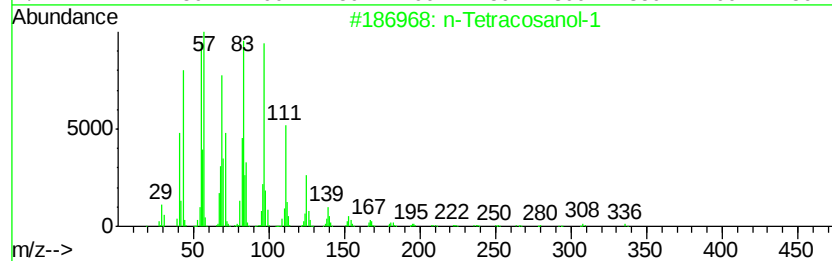
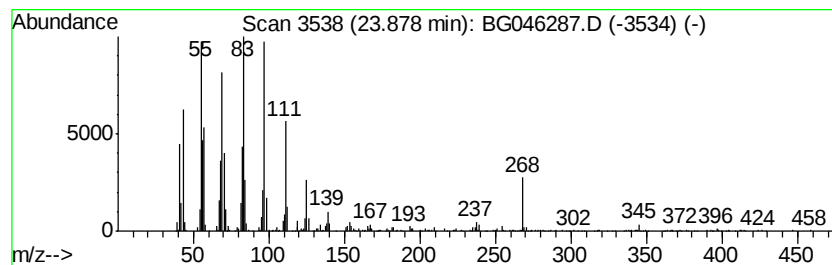
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Tetracosanol-1 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.88	15.15 ng	1022100	Perylene-d12	24.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
2			Behenic alcohol	326	C22H46O	000661-19-8	91
3			1-Heptacosanol	396	C27H56O	002004-39-9	91
4			1-Heneicosanol	312	C21H44O	015594-90-8	91
5			1-Octadecanol	270	C18H38O	000112-92-5	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081420\
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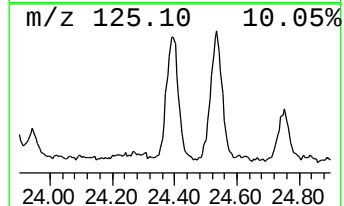
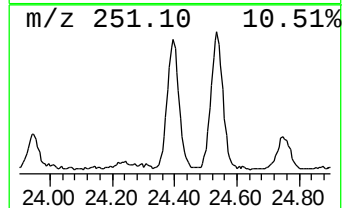
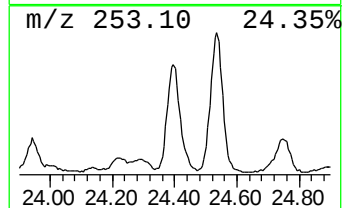
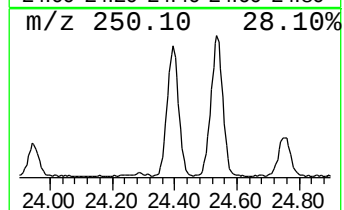
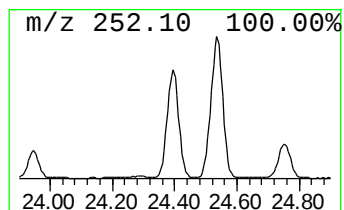
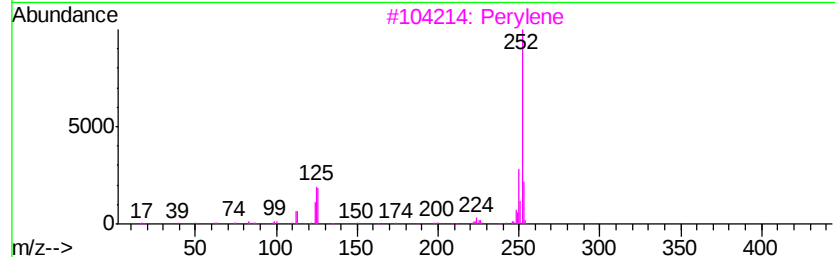
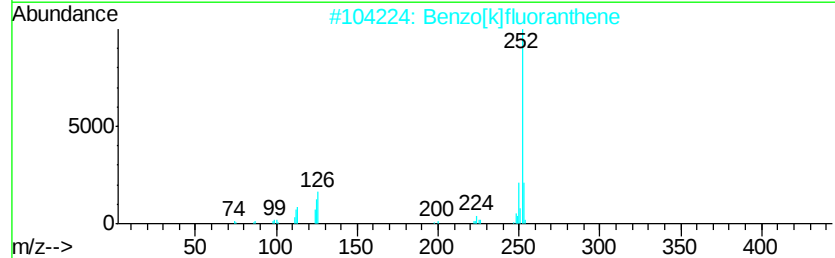
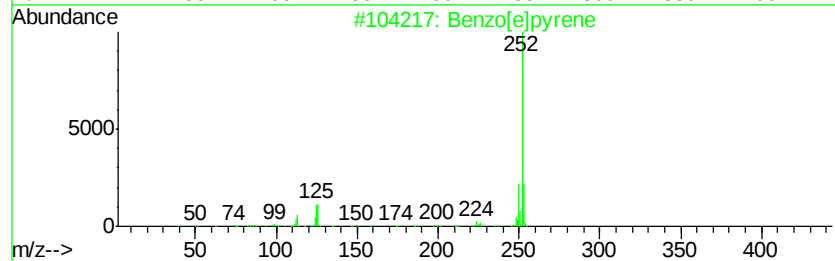
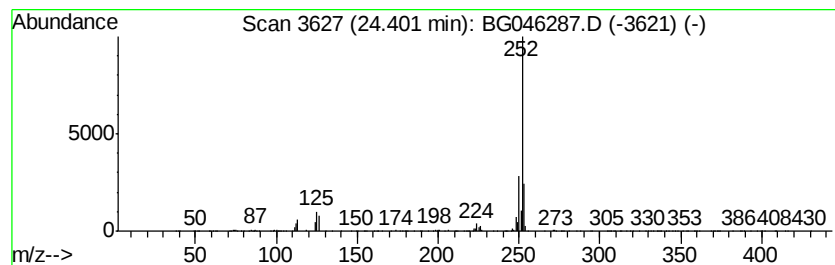
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.40	19.56 ng	1319830	Perylene-d12	24.68

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	97
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
3	Perylene	252	C20H12	000198-55-0	93
4	Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	93
5	Benzo[a]pyrene	252	C20H12	000050-32-8	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081420\
Data File : BG046287.D
Acq On : 14 Aug 2020 22:43
Operator : CG/JU
Sample : L3662-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
VB16182

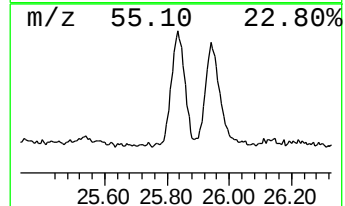
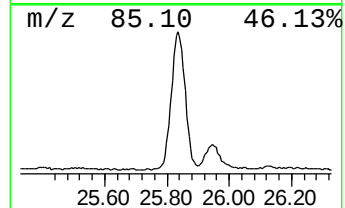
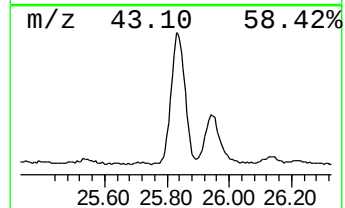
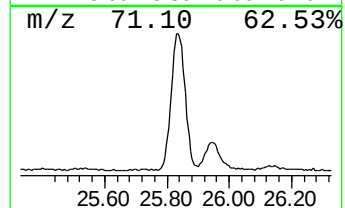
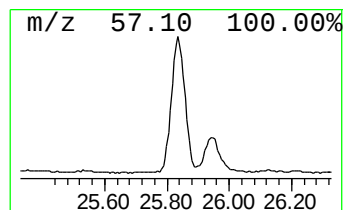
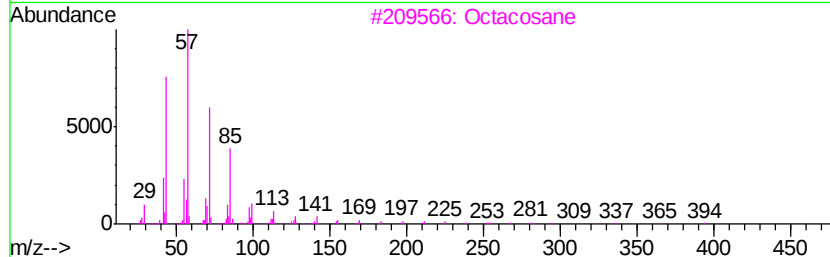
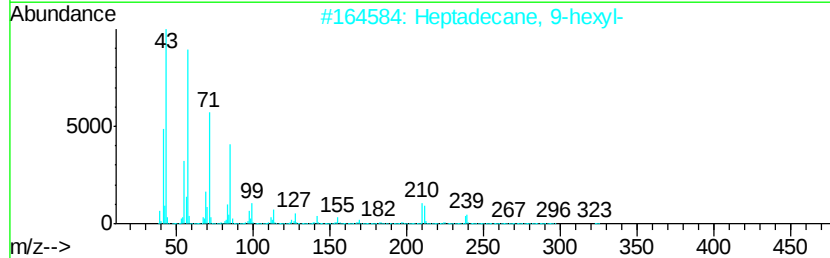
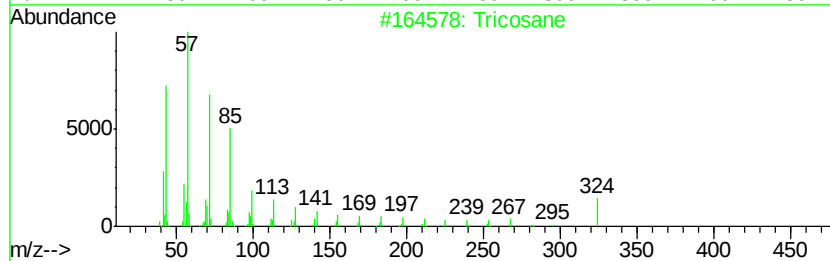
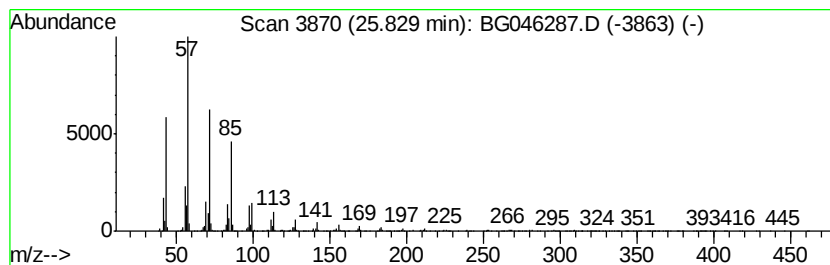
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptadecane, 9-hexyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.83	32.44 ng	2188750	Perylene-d12	24.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosane	324	C23H48	000638-67-5	95
2		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
3		Octacosane	394	C28H58	000630-02-4	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Docosane, 7-hexyl-	394	C28H58	055373-86-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081420\
Data File : BG046287.D
Acq On : 14 Aug 2020 22:43
Operator : CG/JU
Sample : L3662-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

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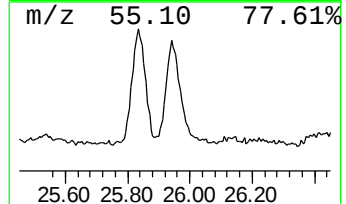
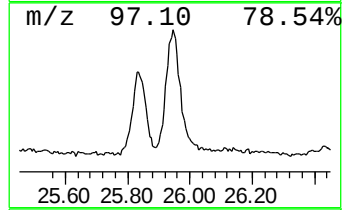
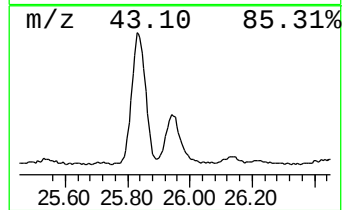
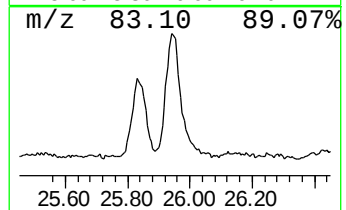
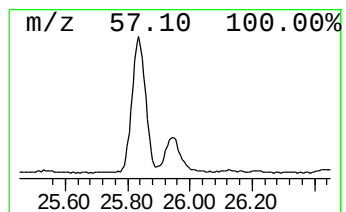
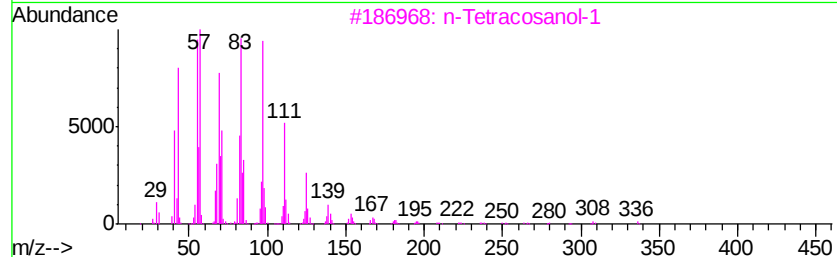
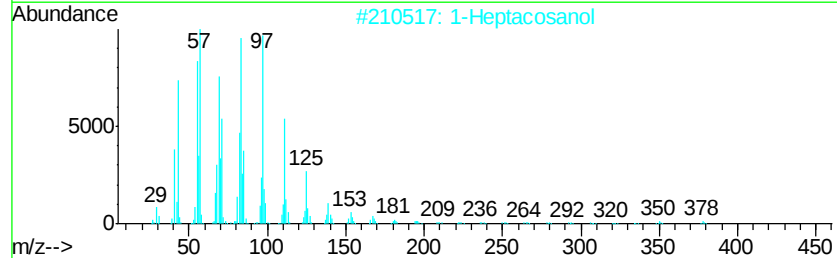
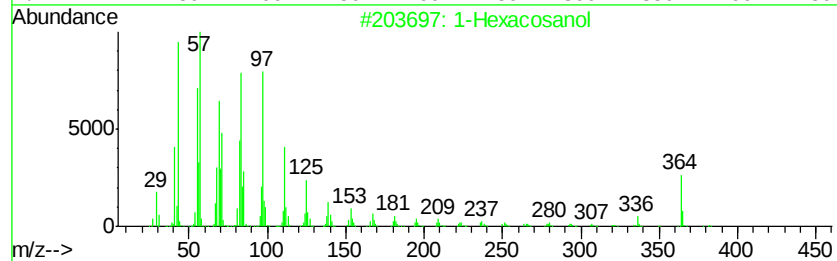
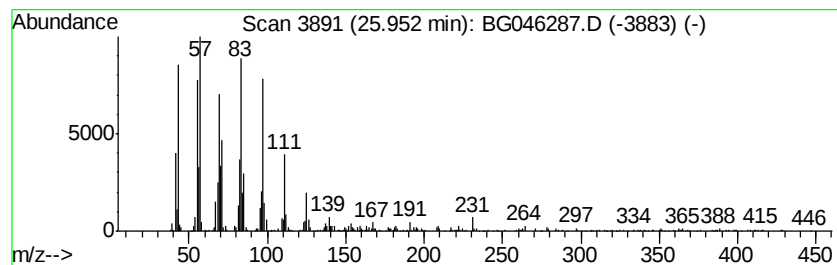
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1-Hexacosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.95	22.91 ng	1545620	Perylene-d12	24.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexacosanol	382	C26H54O	000506-52-5	96
2		1-Heptacosanol	396	C27H56O	002004-39-9	95
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4		Octacosanol	410	C28H58O	000557-61-9	94
5		1-Heneicosanol	312	C21H44O	015594-90-8	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081420\
Data File : BG046287.D
Acq On : 14 Aug 2020 22:43
Operator : CG/JU
Sample : L3662-02
Misc :
ALS Vial : 14 Sample Multiplier: 1

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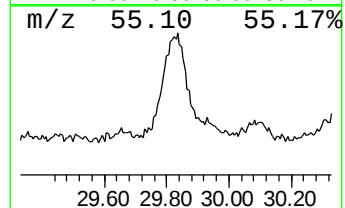
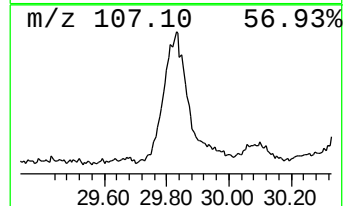
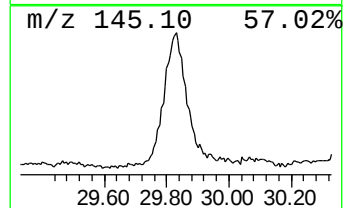
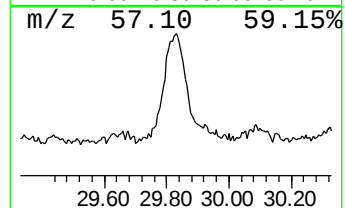
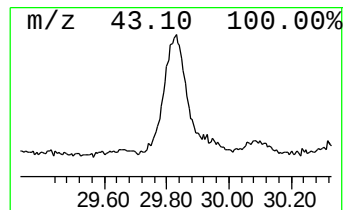
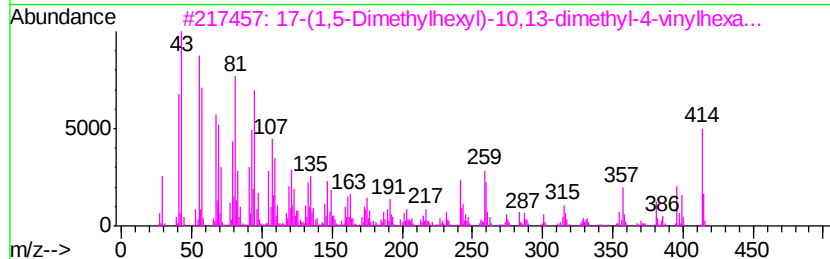
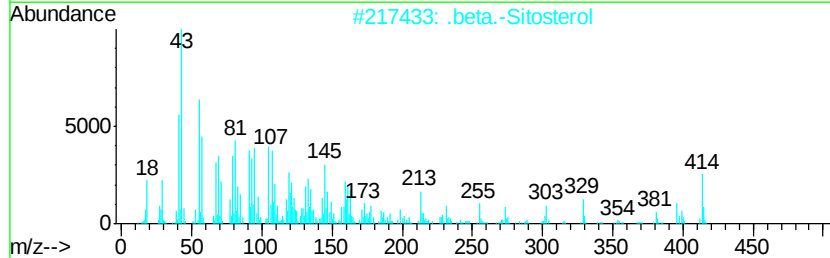
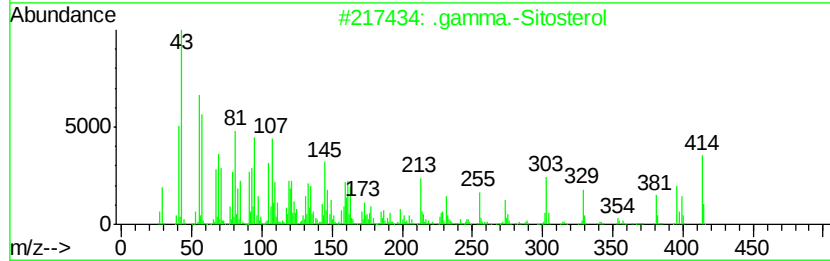
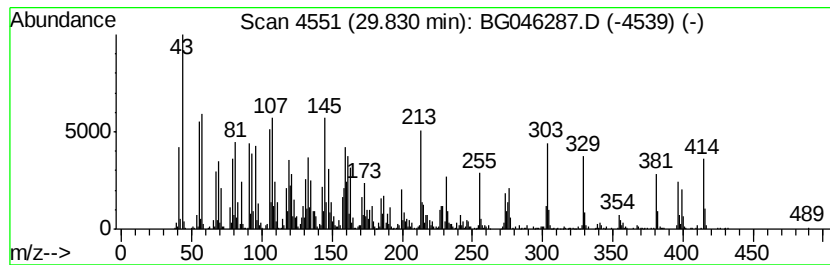
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 .gamma.-Sitosterol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.83	50.68 ng	3418970	Perylene-d12	24.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.gamma.-Sitosterol	414	C29H50O	000083-47-6	98
2		.beta.-Sitosterol	414	C29H50O	000083-46-5	51
3		17-(1,5-Dimethylhexyl)-10,13-dim...	414	C29H50O	1000210-86-9	15
4		Diethylbis(tetramethylene)disilo...	242	C12H26OSi2	018035-51-3	10
5		5.alpha.-Cholest-8-en-3-one, 6.a...	456	C30H48O3	005259-17-6	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG081420\
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 Sample : L3662-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG073020.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
Butane, 2-methoxy...	3.16	14.4	ng	478822	1	7.95	666077	20.0
Cyclohexane, 1,1,...	11.17	2.8	ng	145402	2	10.75	1050830	20.0
unknown12.49	12.49	3.6	ng	189414	2	10.75	1050830	20.0
Decahydro-4,4,8,9...	13.38	4.0	ng	633289	3	14.58	3152550	20.0
Benzene, 1-cycloh...	13.86	2.8	ng	435287	3	14.58	3152550	20.0
2-Pentanone, 4-cy...	14.42	9.9	ng	1568310	3	14.58	3152550	20.0
2,6-Naphthalenedi...	15.38	15.8	ng	2483770	3	14.58	3152550	20.0
Octadecane	22.38	7.0	ng	1031260	5	21.58	2948490	20.0
Strychnidin-10-on...	22.44	20.2	ng	2983180	5	21.58	2948490	20.0
Tricosane	23.83	26.7	ng	1801380	6	24.68	1349260	20.0
n-Tetracosanol-1	23.88	15.2	ng	1022100	6	24.68	1349260	20.0
Benzo[e]pyrene	24.40	19.6	ng	1319830	6	24.68	1349260	20.0
Heptadecane, 9-he...	25.83	32.4	ng	2188750	6	24.68	1349260	20.0
1-Hexacosanol	25.95	22.9	ng	1545620	6	24.68	1349260	20.0
.gamma.-Sitosterol	29.83	50.7	ng	3418970	6	24.68	1349260	20.0