

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070721\
 Data File : BG049198.D
 Acq On : 7 Jul 2021 14:00
 Operator : CG/JU
 Sample : M2933-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 3

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-BG063021.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG049198.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.128	9	15	26	rVV3	53630	161881	9.70%	1.324%
2	3.210	26	29	36	rVB6	17634	36357	2.18%	0.297%
3	5.173	356	363	369	rBV	12437	21816	1.31%	0.178%
4	5.643	436	443	453	rBV	434898	737735	44.20%	6.033%
5	7.241	706	715	725	rBV	456299	797115	47.75%	6.518%
6	8.087	851	859	866	rBV2	133794	234167	14.03%	1.915%
7	9.250	1050	1057	1065	rBV	281008	506914	30.37%	4.145%
8	10.907	1331	1339	1348	rVB	171512	315043	18.87%	2.576%
9	13.351	1746	1755	1762	rBV	722494	1110243	66.51%	9.079%
10	14.174	1888	1895	1901	rBV	77932	120515	7.22%	0.985%
11	14.726	1983	1989	1997	rVB	253924	422243	25.30%	3.453%
12	16.219	2236	2243	2252	rBV	663597	1019894	61.10%	8.340%
13	17.476	2451	2457	2464	rVB	326330	523539	31.36%	4.281%
14	18.357	2597	2607	2617	rVB2	106816	190229	11.40%	1.556%
15	18.704	2657	2666	2676	rVB	68258	255816	15.33%	2.092%
16	18.839	2680	2689	2701	rBV3	100363	429785	25.75%	3.514%
17	18.957	2701	2709	2721	rVB	291487	714827	42.82%	5.845%
18	19.127	2733	2738	2745	rBV3	22712	47967	2.87%	0.392%
19	19.327	2769	2772	2779	rVB6	16737	28625	1.71%	0.234%
20	19.620	2819	2822	2835	rVB2	157865	280346	16.79%	2.292%
21	19.885	2860	2867	2876	rBV2	63092	137675	8.25%	1.126%
22	20.085	2895	2901	2906	rVB	1304609	1669271	100.00%	13.650%
23	20.690	2998	3004	3013	rBV	86202	173288	10.38%	1.417%
24	21.401	3120	3125	3133	rBV3	109488	239102	14.32%	1.955%
25	21.483	3133	3139	3148	rVB2	108984	225822	13.53%	1.847%
26	21.759	3180	3186	3193	rVB	344102	596186	35.72%	4.875%
27	22.382	3286	3292	3300	rVB	80914	179060	10.73%	1.464%
28	22.582	3322	3326	3329	rBV5	11659	21528	1.29%	0.176%
29	23.422	3462	3469	3477	rVB2	55822	141117	8.45%	1.154%
30	24.121	3584	3588	3592	rBV7	17340	30481	1.83%	0.249%
31	24.620	3666	3673	3674	rBV6	40197	75938	4.55%	0.621%
32	25.067	3738	3749	3764	rVB3	209876	689322	41.29%	5.637%
33	26.072	3912	3920	3921	rBV4	38499	70186	4.20%	0.574%
34	27.787	4208	4212	4213	rBV4	18496	25080	1.50%	0.205%

Sum of corrected areas: 12229113

Instrument :

BNA_G

ClientSampleId :

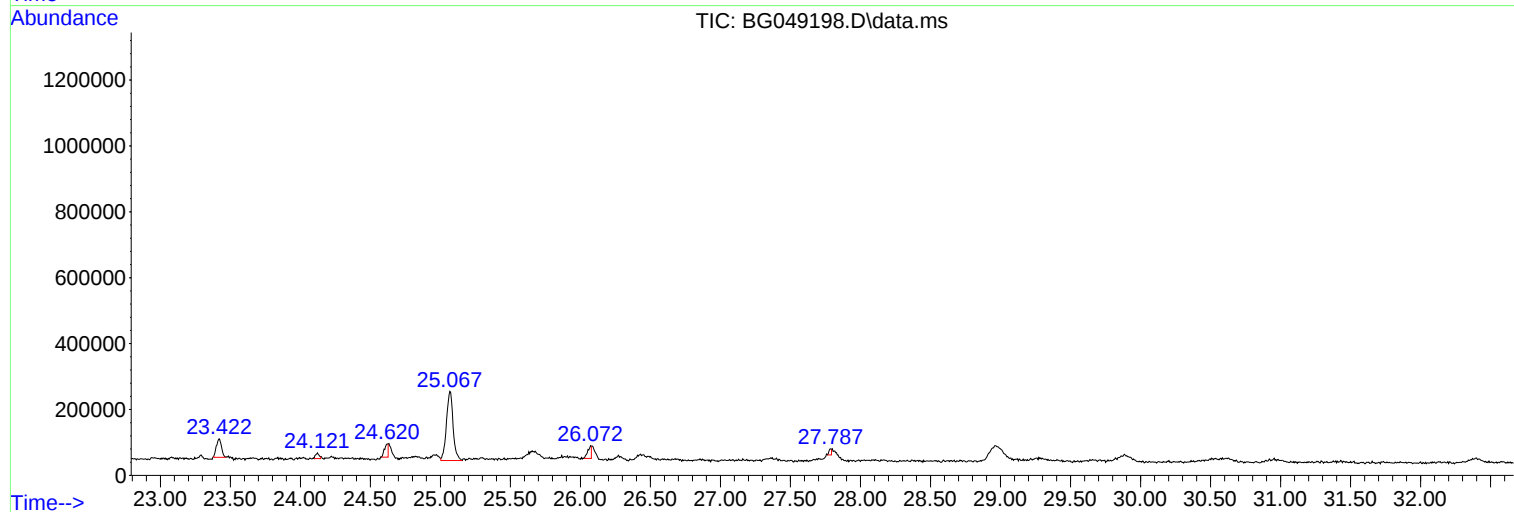
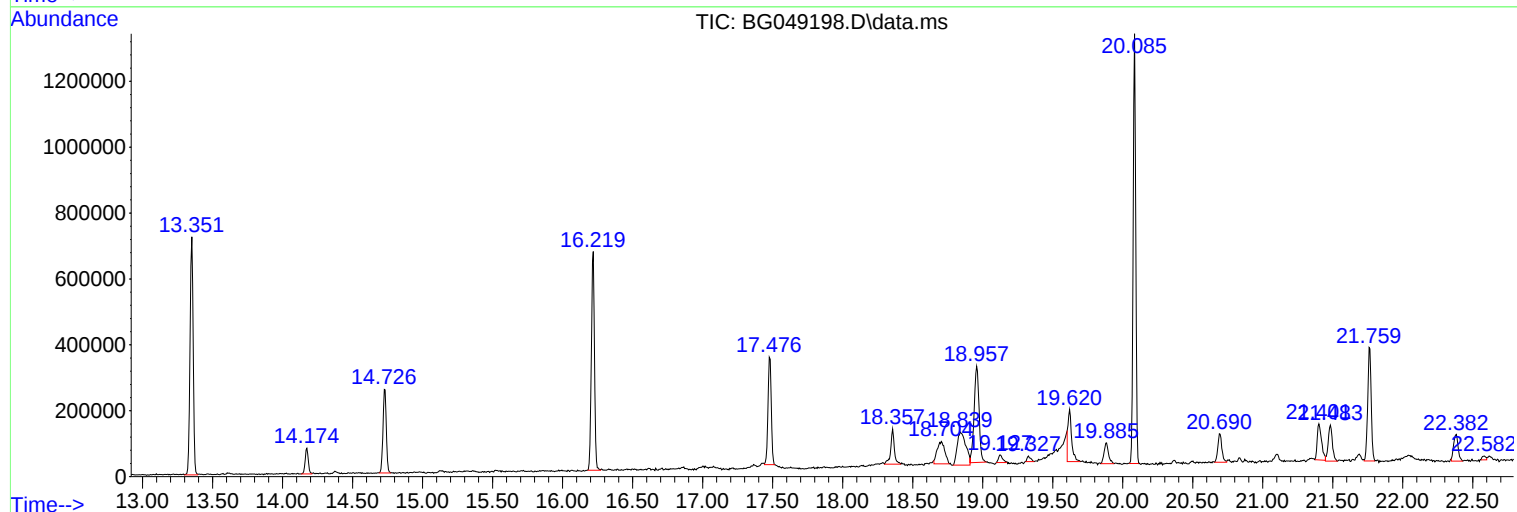
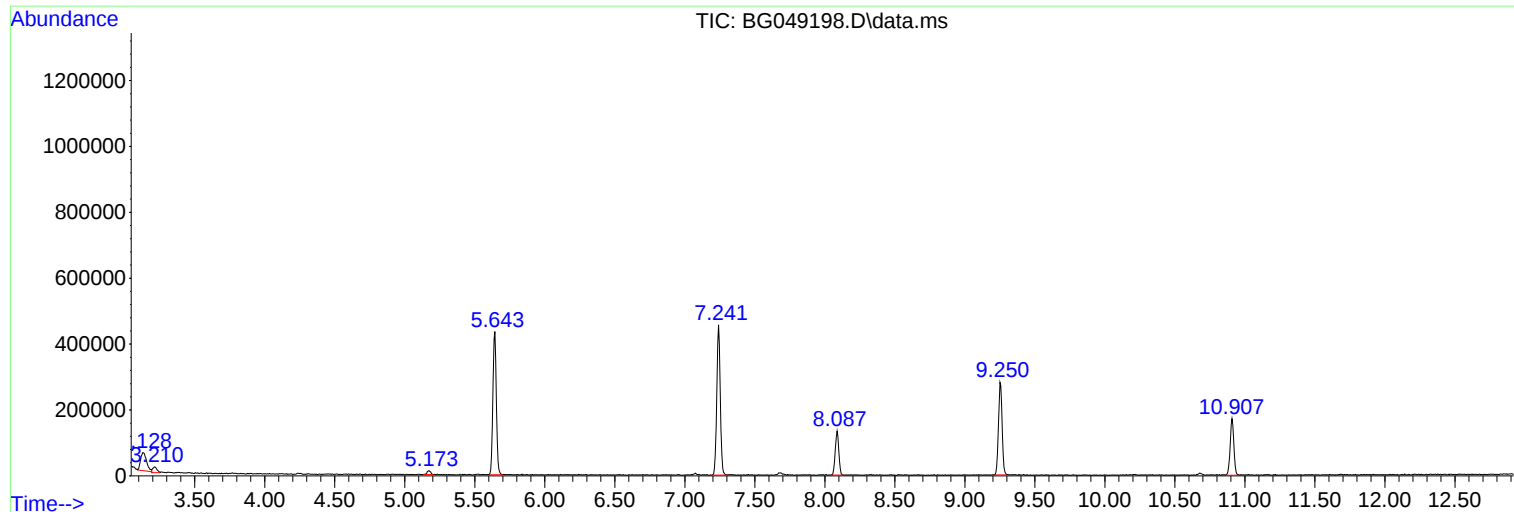
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG063021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Operator : CG/JU
Sample : M2933-03
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Instrument :
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ClientSampled :
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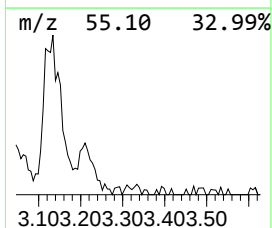
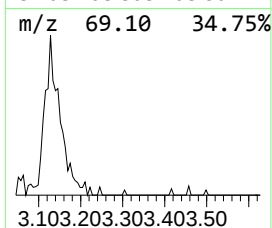
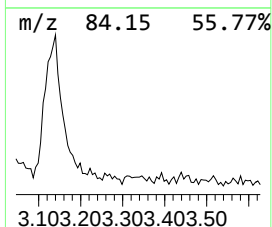
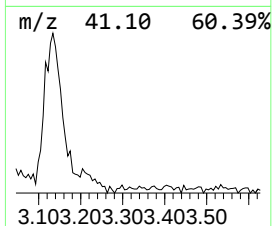
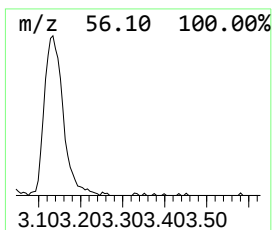
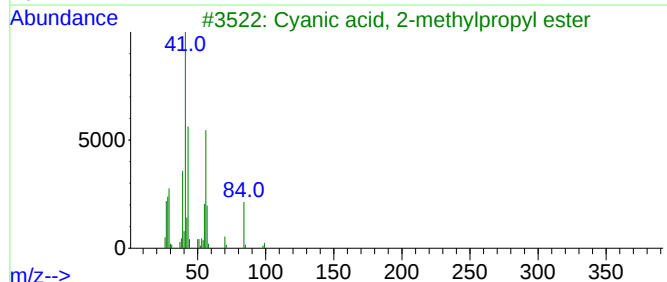
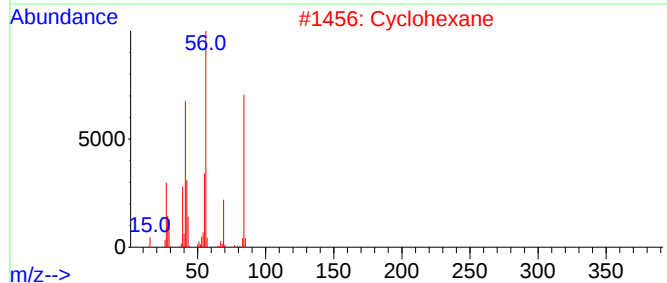
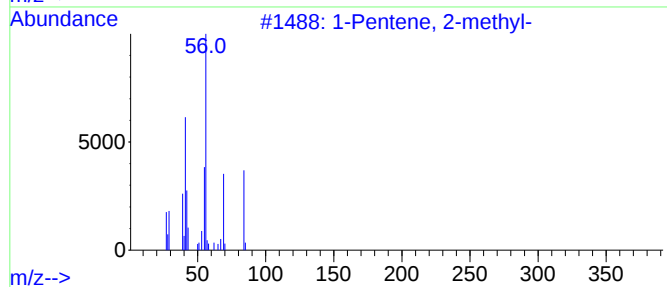
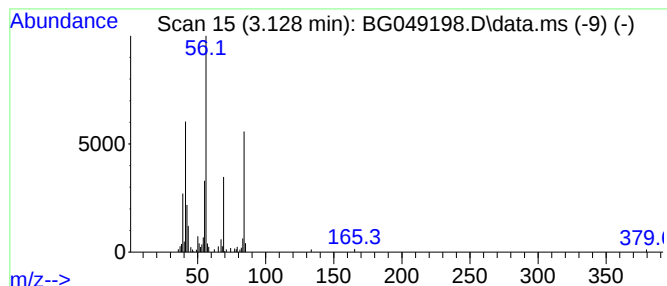
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 1-Pentene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.128	13.83 ng	161881	1,4-Dichlorobenzene-d4	8.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	90
2			Cyclohexane	84	C6H12	000110-82-7	90
3			Cyanic acid, 2-methylpropyl ester	99	C5H9NO	001768-25-8	64
4			Cyclobutane, ethyl-	84	C6H12	004806-61-5	43
5			Cyclopentane, methyl-	84	C6H12	000096-37-7	42



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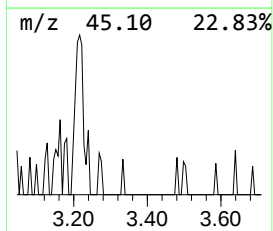
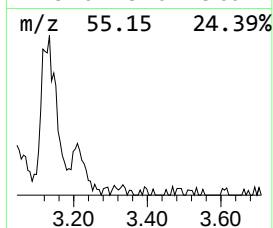
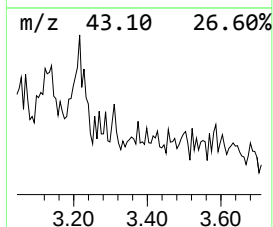
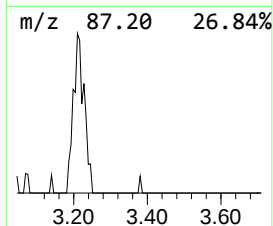
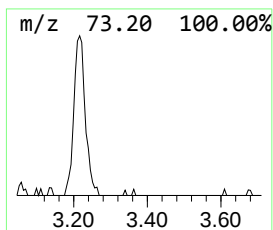
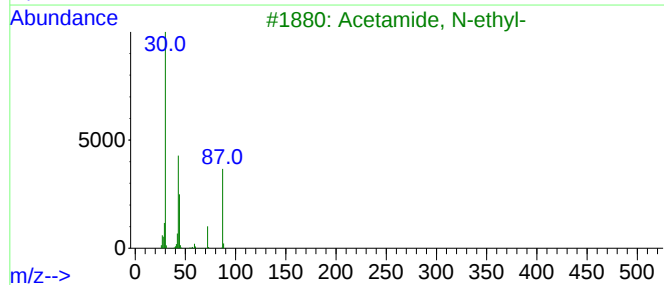
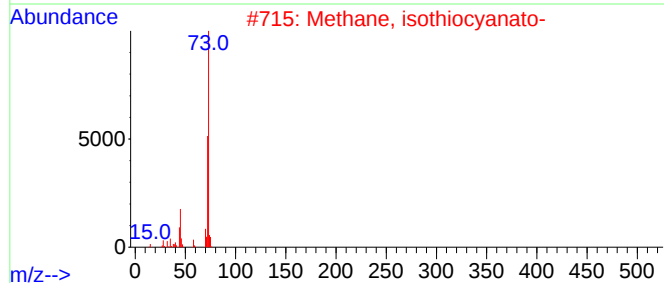
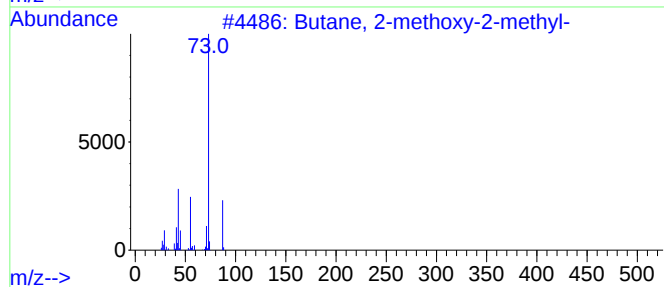
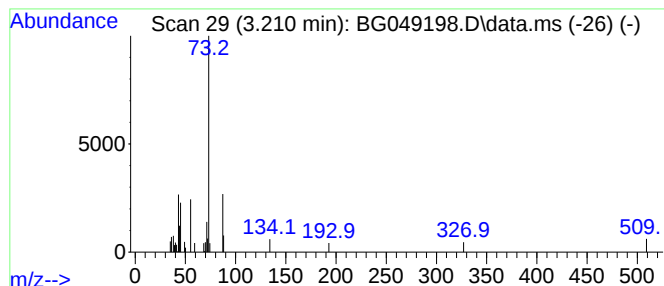
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG063021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
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Peak Number 2 unknown3.210 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.210	3.11 ng	36357	1,4-Dichlorobenzene-d4	8.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	38
2			Methane, isothiocyanato-	73	C2H3NS	000556-61-6	9
3			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
4			1,5-Pentanediamine	102	C5H14N2	000462-94-2	9
5			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	9



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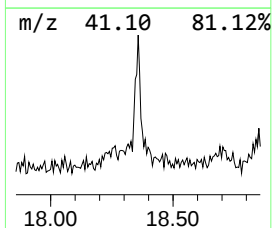
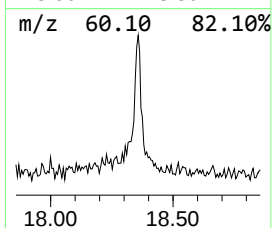
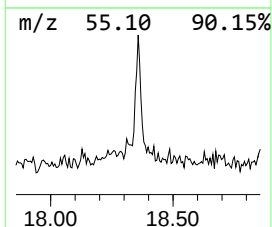
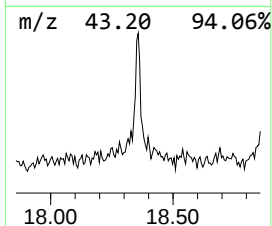
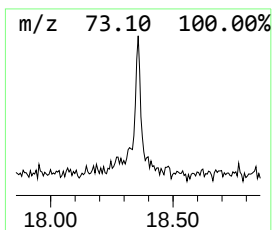
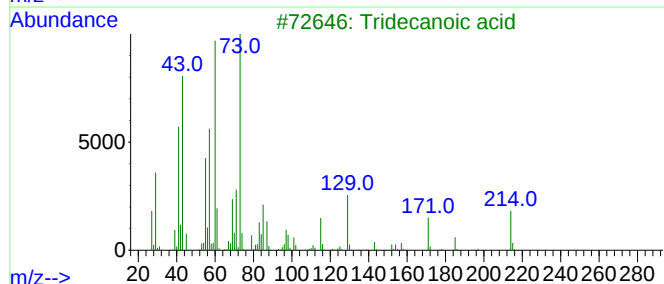
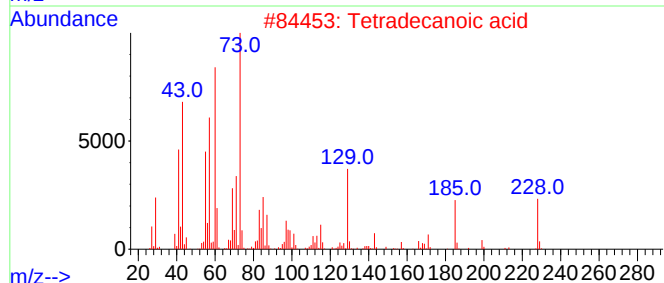
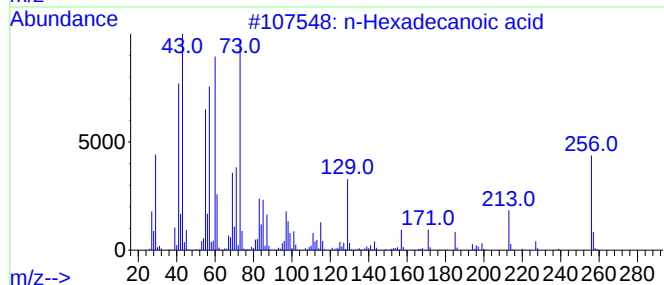
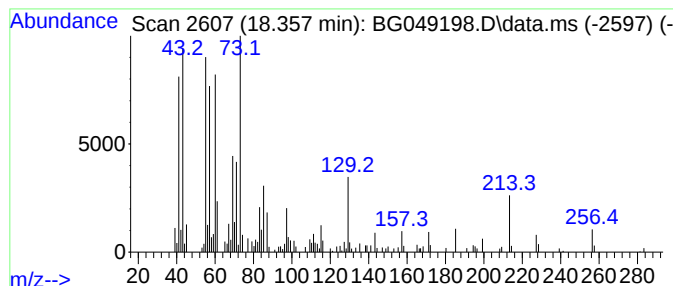
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.357	7.27 ng	190229	Phenanthrene-d10	17.482

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	81
3			Tridecanoic acid	214	C13H26O2	000638-53-9	64
4			n-Decanoic acid	172	C10H20O2	000334-48-5	50
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	35



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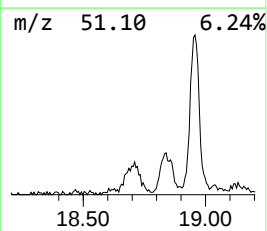
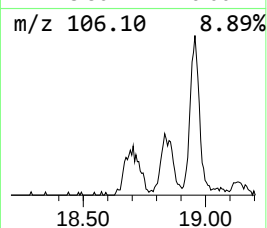
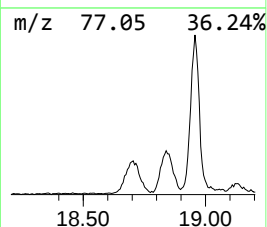
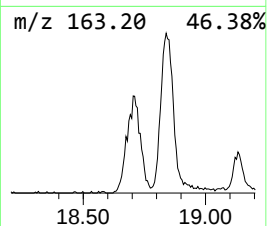
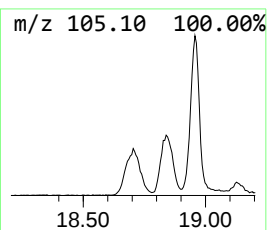
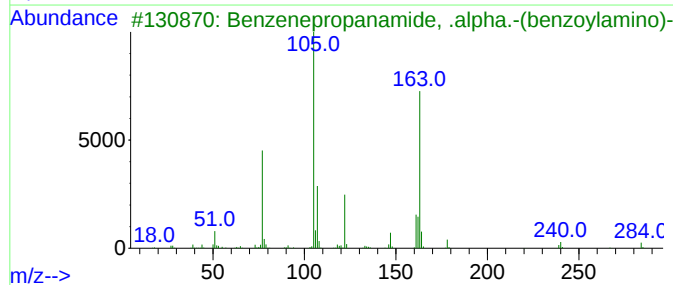
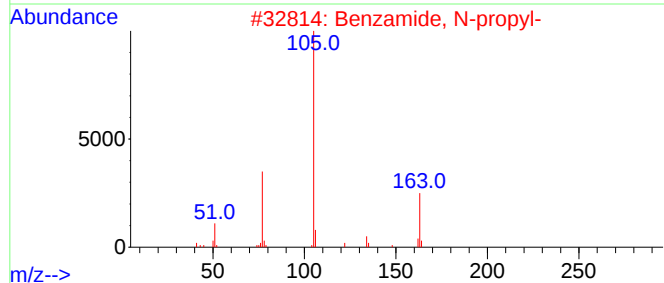
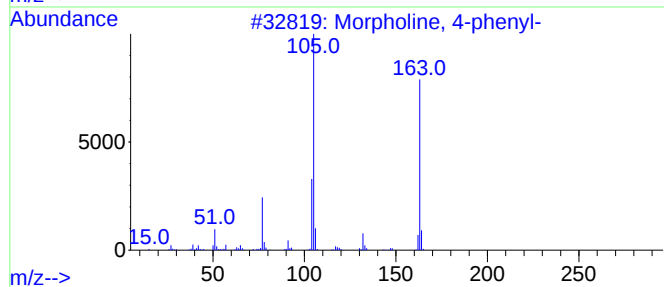
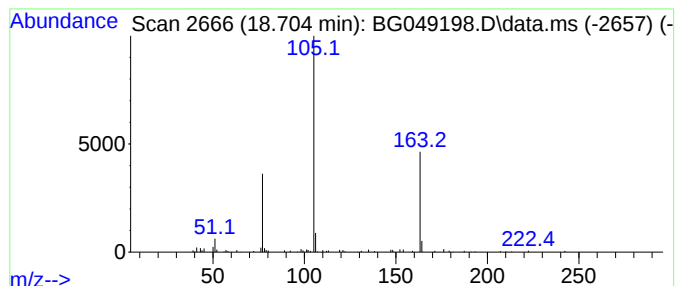
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 Morpholine, 4-phenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.704	9.77 ng	255816	Phenanthrene-d10	17.482

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	56
2		Benzamide, N-propyl-	163	C10H13NO	010546-70-0	56
3		Benzenepropanamide, .alpha.-(ben...	284	C16H16N2O3	058690-81-6	56
4		1-Butanone, 1-phenyl-	148	C10H12O	000495-40-9	27
5		N-(m-Nitrophenyl)benzamide	242	C13H10N2O3	004771-08-8	27



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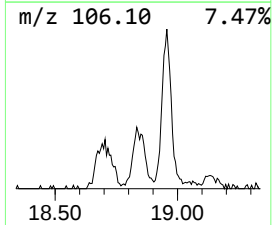
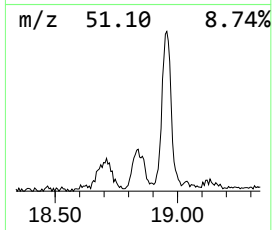
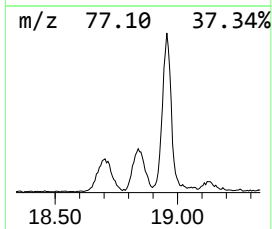
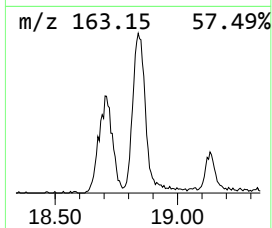
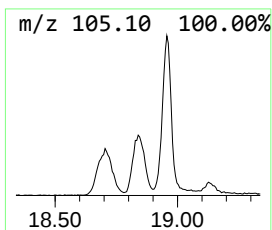
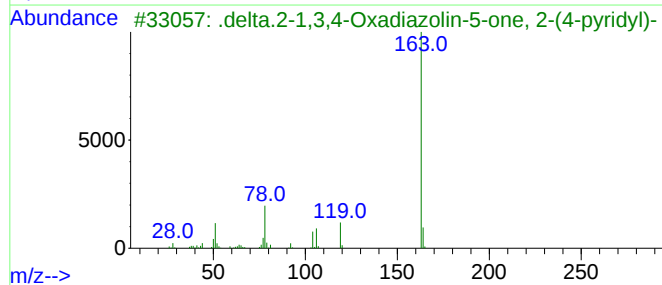
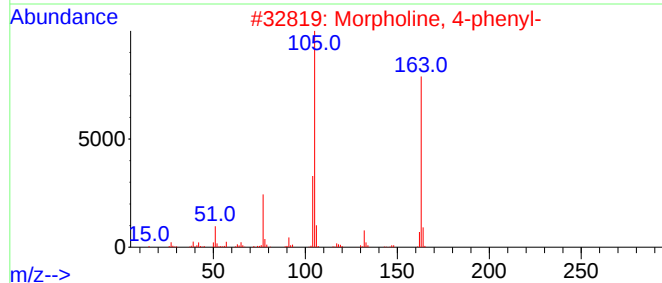
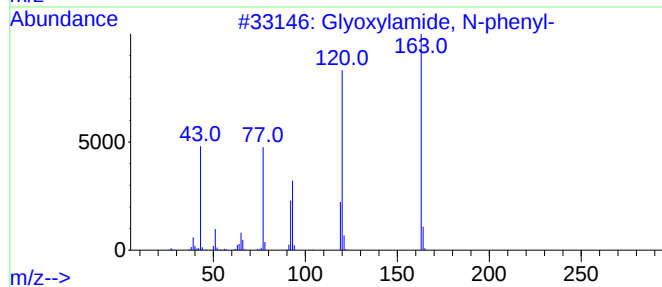
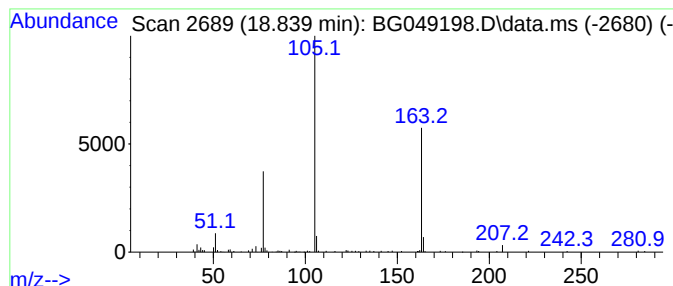
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TIC Library : C:\Database\NIST11.L
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Peak Number 5 unknown18.839 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.839	16.42 ng	429785	Phenanthrene-d10	17.482

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Glyoxylamide, N-phenyl-	163	C9H9NO2	032331-52-5	47
2			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	40
3			.delta.2-1,3,4-Oxadiazolin-5-one...	163	C7H5N3O2	002845-82-1	38
4			1,3-Propanediol dibenzoate	284	C17H16O4	002451-86-7	35
5			4'-Nitrobenzanilide	242	C13H10N2O3	003393-96-2	27



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ClientSampled :
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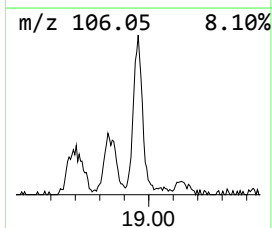
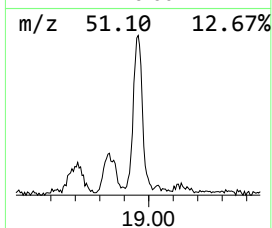
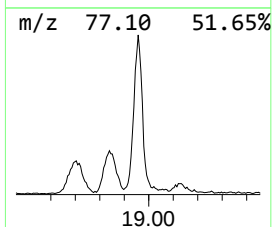
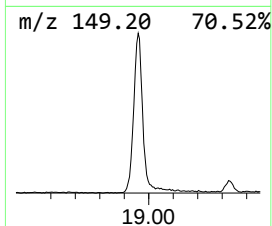
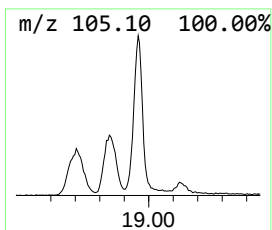
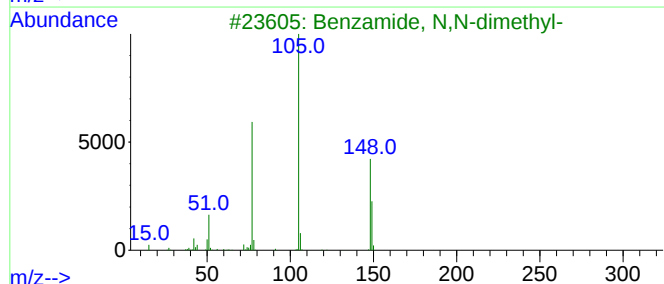
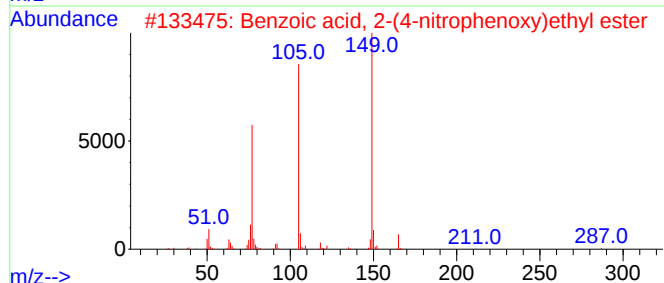
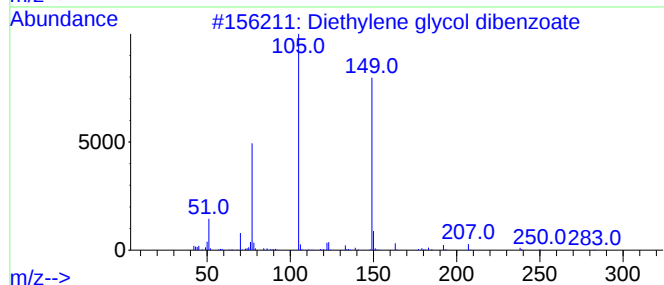
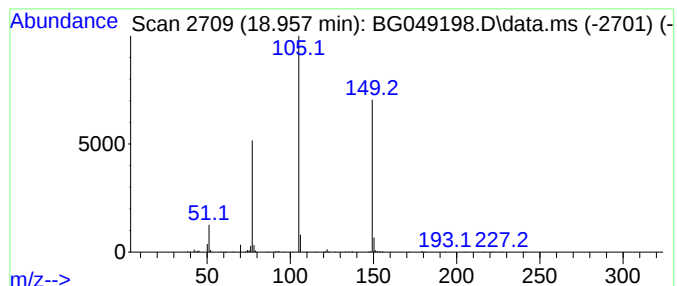
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Diethylene glycol dibenzoate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.956	27.31 ng	714827	Phenanthrene-d10	17.482

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	83
2			Benzoic acid, 2-(4-nitrophenoxy)...	287	C15H13NO5	1000368-92-7	72
3			Benzamide, N,N-dimethyl-	149	C9H11NO	000611-74-5	53
4			1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	50
5			Ethanone, 2,2,2-trifluoro-1-phenyl-	174	C8H5F3O	000434-45-7	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070721\
Data File : BG049198.D
Acq On : 7 Jul 2021 14:00
Operator : CG/JU
Sample : M2933-03
Misc :
ALS Vial : 5 Sample Multiplier: 1

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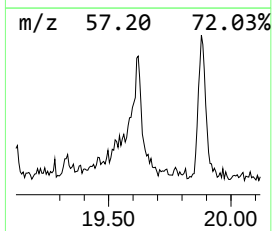
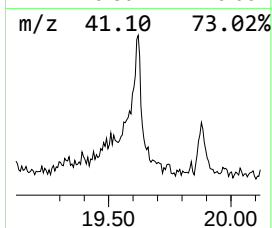
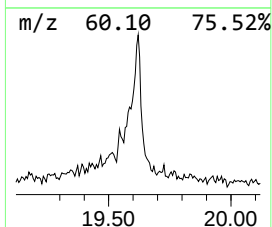
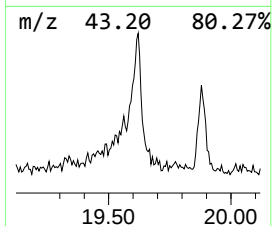
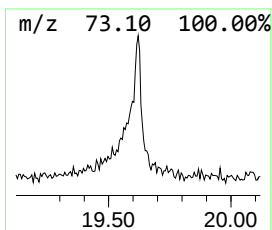
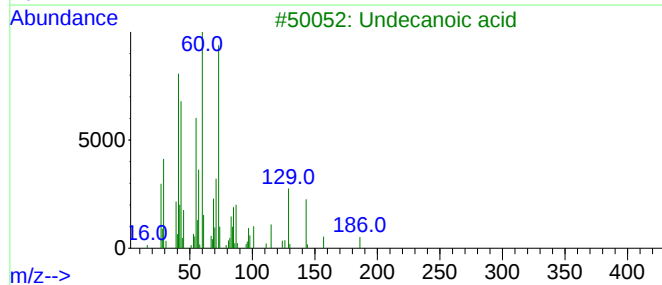
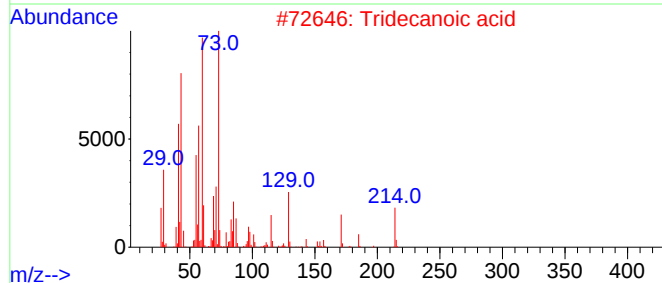
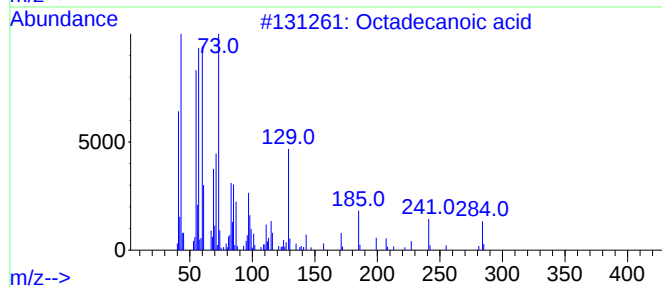
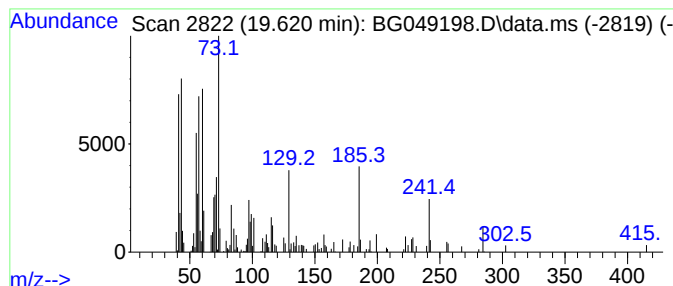
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.620	10.71 ng	280346	Phenanthrene-d10	17.482

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	89
2			Tridecanoic acid	214	C13H26O2	000638-53-9	50
3			Undecanoic acid	186	C11H22O2	000112-37-8	46
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	42
5			Undecanoic acid, 10-bromo-	264	C11H21BrO2	018294-93-4	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070721\
Data File : BG049198.D
Acq On : 7 Jul 2021 14:00
Operator : CG/JU
Sample : M2933-03
Misc :
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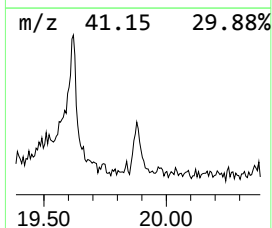
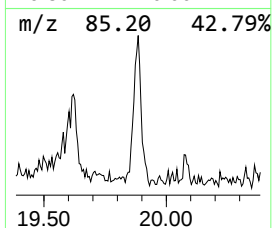
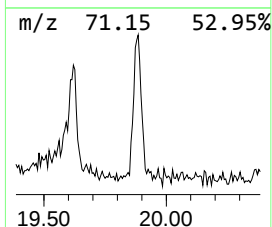
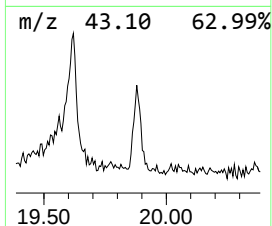
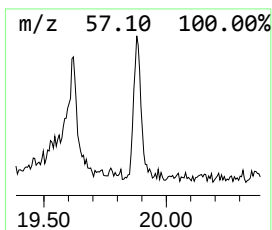
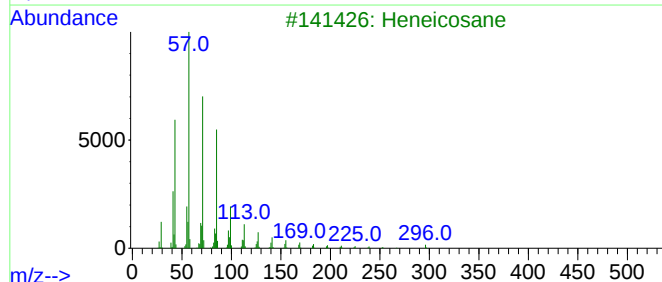
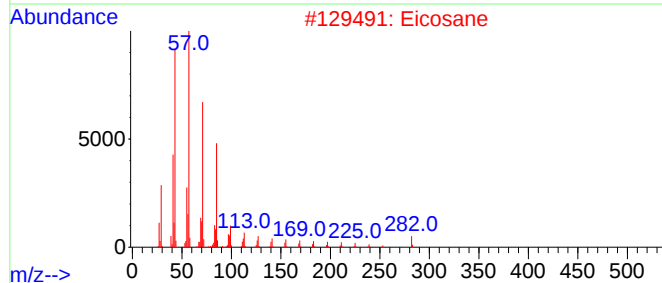
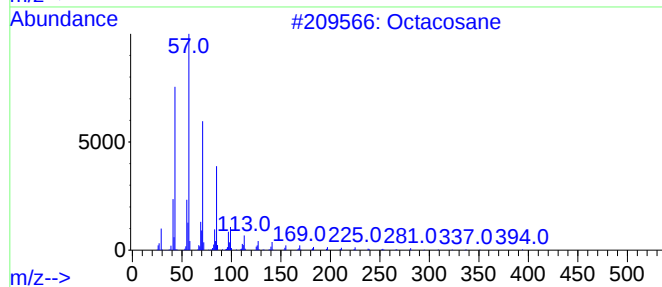
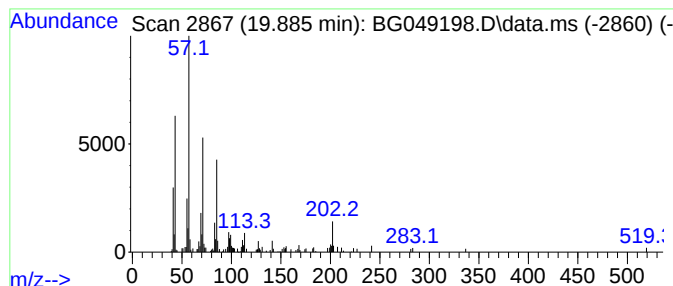
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG063021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.885	4.62 ng	137675	Chrysene-d12	21.759	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	81
2	Eicosane	282	C20H42	000112-95-8	81
3	Heneicosane	296	C21H44	000629-94-7	74
4	Tetracosane	338	C24H50	000646-31-1	74
5	Heptadecane	240	C17H36	000629-78-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070721\
Data File : BG049198.D
Acq On : 7 Jul 2021 14:00
Operator : CG/JU
Sample : M2933-03
Misc :
ALS Vial : 5 Sample Multiplier: 1

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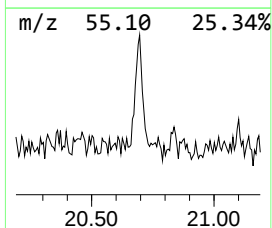
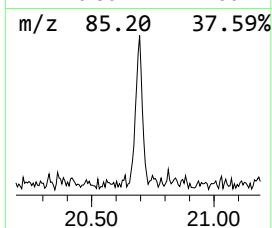
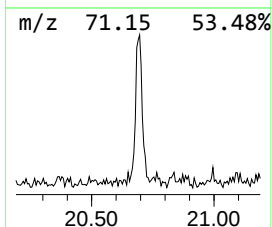
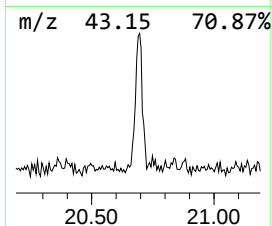
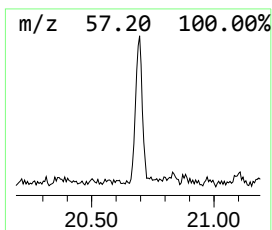
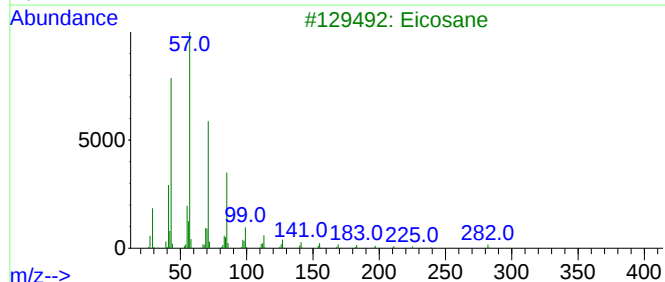
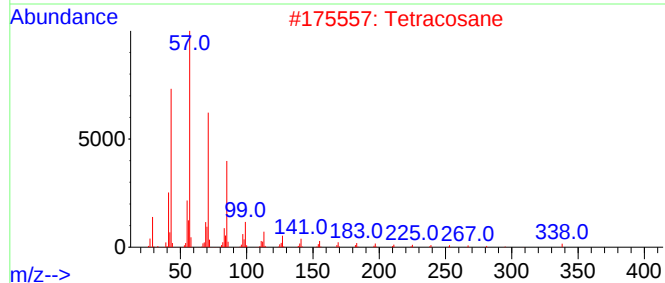
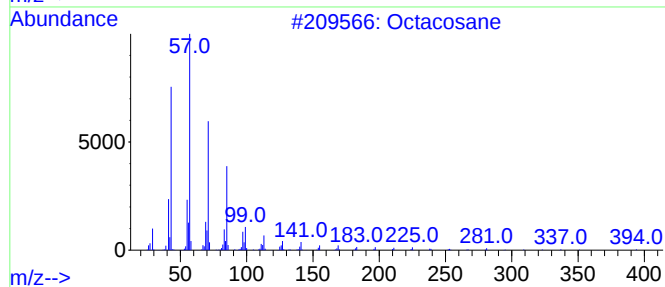
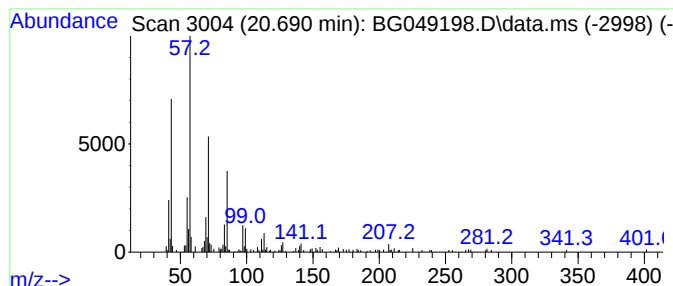
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetracosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.690	5.81 ng	173288	Chrysene-d12	21.759

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	81
2			Tetracosane	338	C24H50	000646-31-1	80
3			Eicosane	282	C20H42	000112-95-8	80
4			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	74
5			Heptacosane	380	C27H56	000593-49-7	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070721\
Data File : BG049198.D
Acq On : 7 Jul 2021 14:00
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Sample : M2933-03
Misc :
ALS Vial : 5 Sample Multiplier: 1

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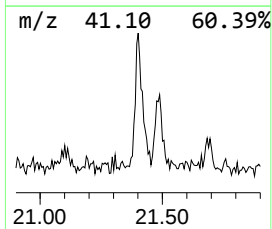
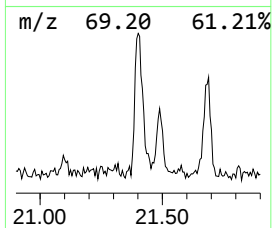
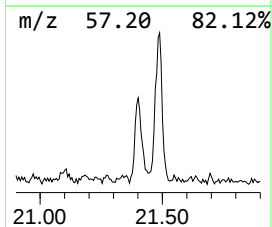
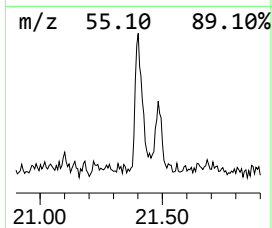
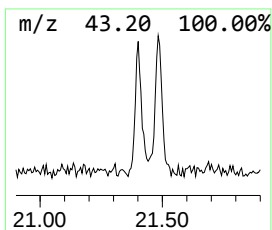
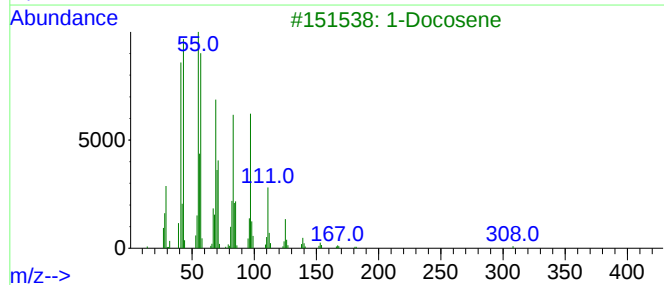
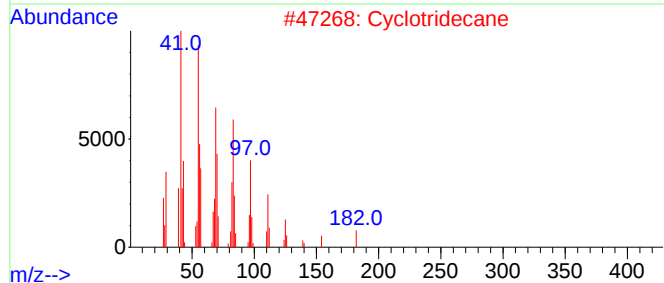
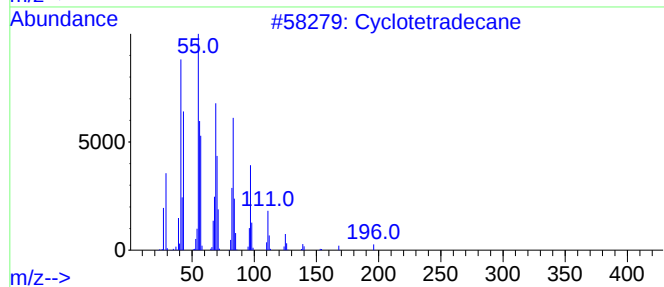
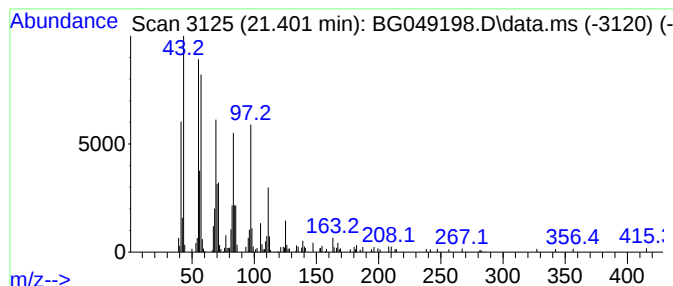
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Cyclotetradecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.401	8.02 ng	239102	Chrysene-d12	21.759

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetradecane	196	C14H28	000295-17-0	96
2			Cyclotridecane	182	C13H26	000295-02-3	94
3			1-Docosene	308	C22H44	001599-67-3	94
4			1-Docosanethiol	342	C22H46S	007773-83-3	91
5			Cyclooctacosane	392	C28H56	000297-24-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070721\
Data File : BG049198.D
Acq On : 7 Jul 2021 14:00
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Sample : M2933-03
Misc :
ALS Vial : 5 Sample Multiplier: 1

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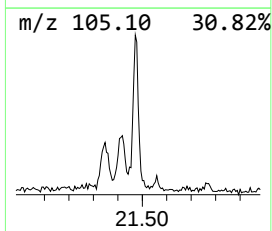
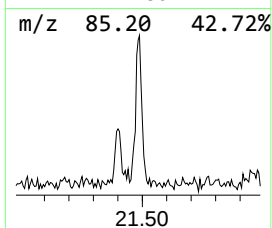
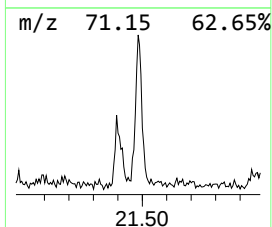
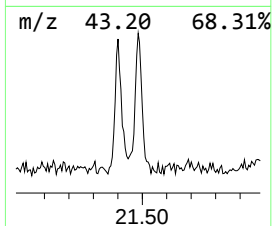
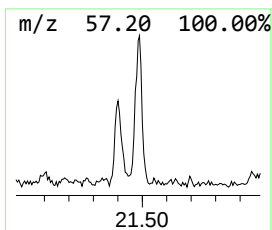
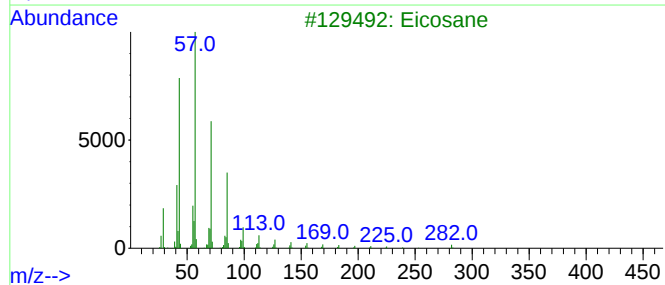
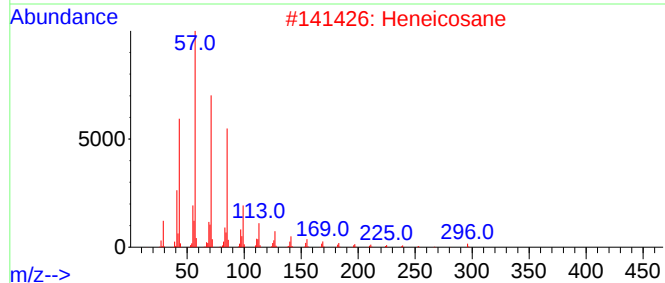
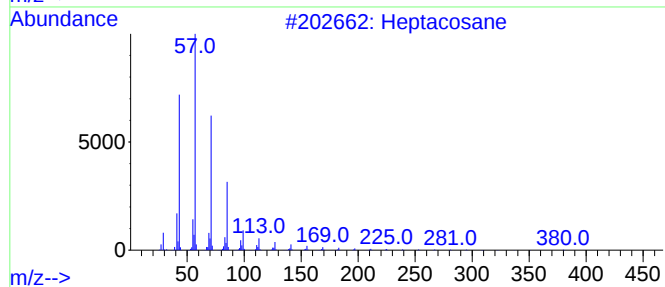
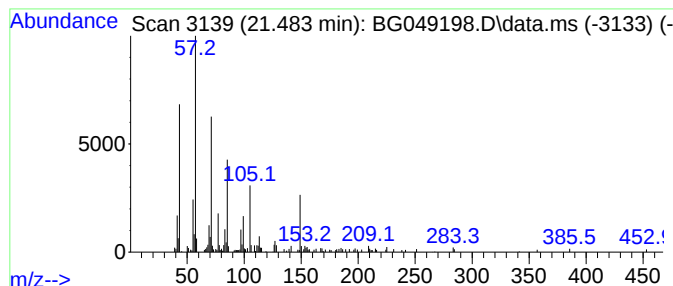
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.483	7.58 ng	225822	Chrysene-d12	21.759

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptacosane	380	C27H56	000593-49-7	58
2		Heneicosane	296	C21H44	000629-94-7	53
3		Eicosane	282	C20H42	000112-95-8	52
4		Nonadecane	268	C19H40	000629-92-5	52
5		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070721\
Data File : BG049198.D
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ALS Vial : 5 Sample Multiplier: 1

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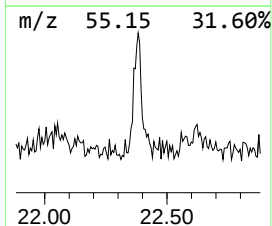
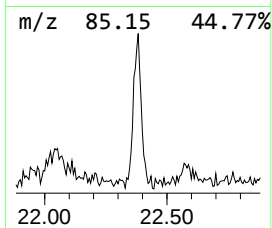
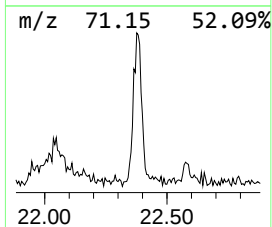
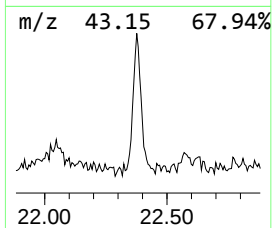
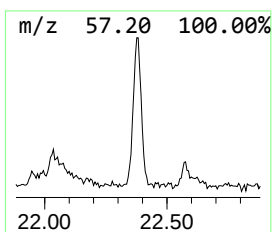
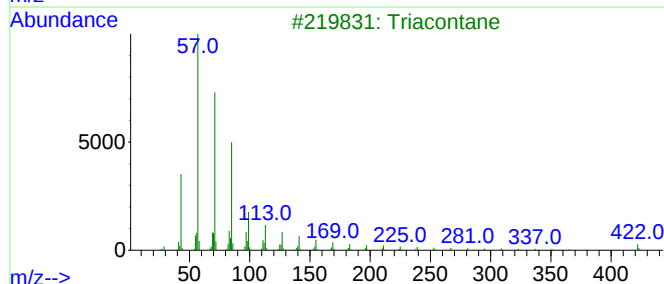
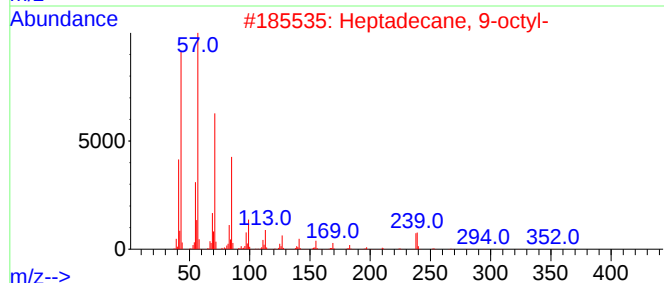
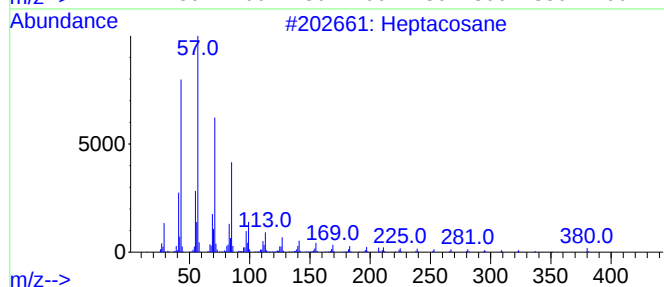
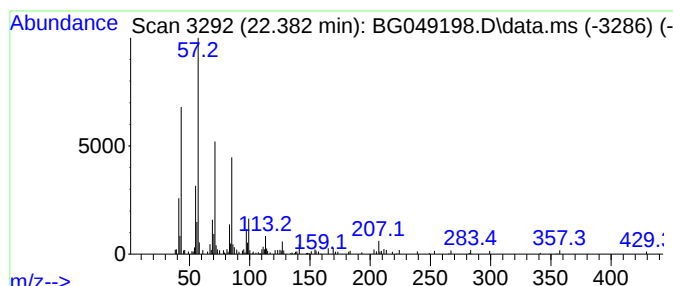
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.382	6.01 ng	179060	Chrysene-d12	21.759

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	87
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	83
3			Triacontane	422	C30H62	000638-68-6	83
4			Heneicosane	296	C21H44	000629-94-7	83
5			2-methyloctacosane	408	C29H60	1000376-72-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070721\
Data File : BG049198.D
Acq On : 7 Jul 2021 14:00
Operator : CG/JU
Sample : M2933-03
Misc :
ALS Vial : 5 Sample Multiplier: 1

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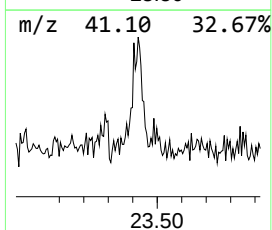
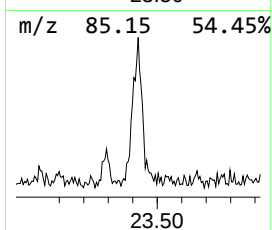
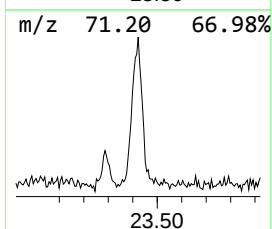
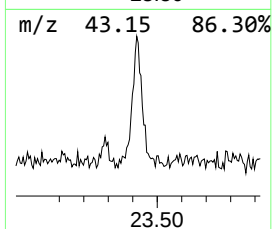
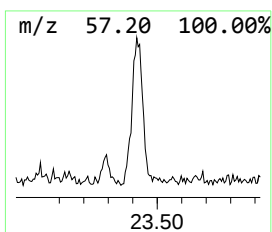
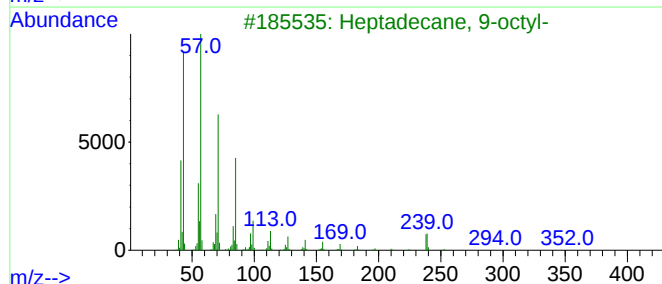
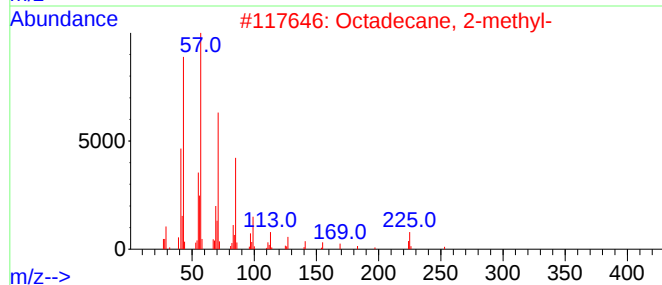
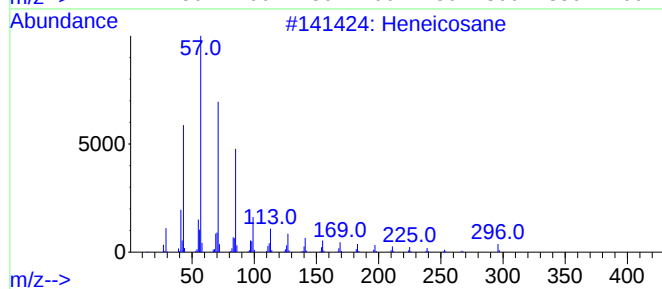
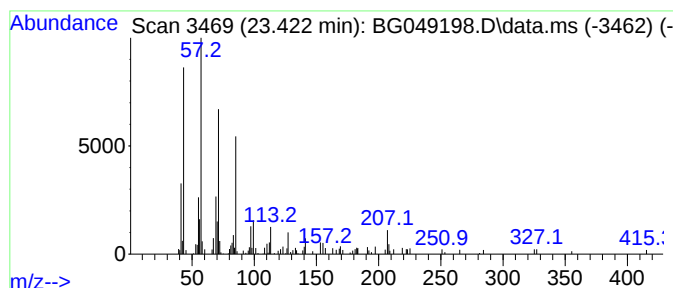
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heneicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.422	4.09 ng	141117	Perylene-d12	25.067

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	83
2		Octadecane, 2-methyl-	268	C19H40	001560-88-9	80
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	80
4		Tetracosane	338	C24H50	000646-31-1	80
5		Tetracosane, 11-decyl-	479	C34H70	055429-84-0	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070721\
Data File : BG049198.D
Acq On : 7 Jul 2021 14:00
Operator : CG/JU
Sample : M2933-03
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
3

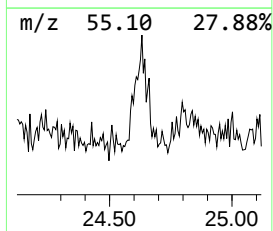
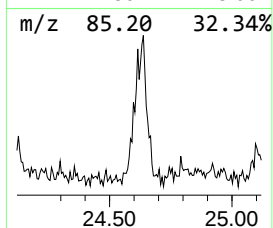
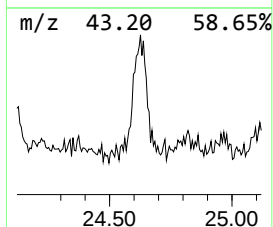
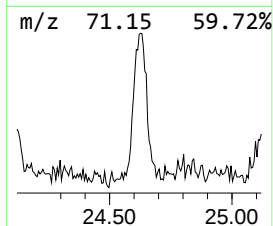
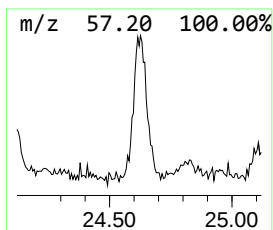
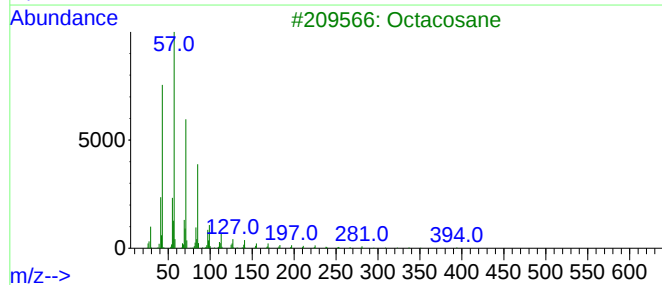
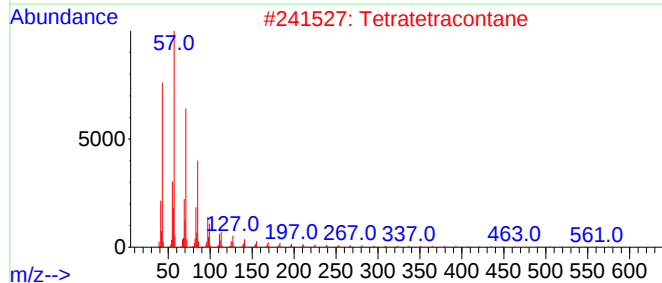
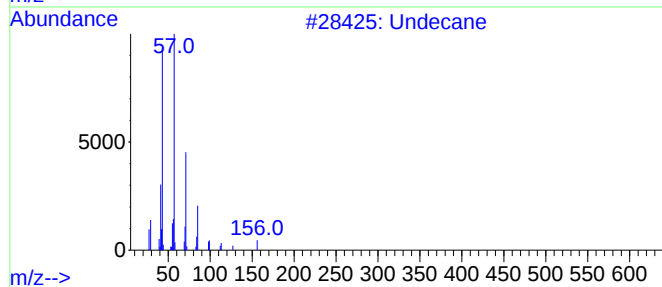
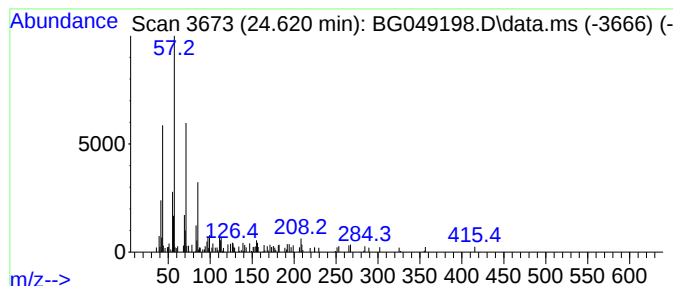
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Undecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.621	2.20 ng	75938	Perylene-d12	25.067

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	68
2	Tetratetracontane			619	C44H90	007098-22-8	58
3	Octacosane			394	C28H58	000630-02-4	58
4	Undecane, 3,8-dimethyl-			184	C13H28	017301-30-3	58
5	Tetradecane			198	C14H30	000629-59-4	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070721\
Data File : BG049198.D
Acq On : 7 Jul 2021 14:00
Operator : CG/JU
Sample : M2933-03
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
3

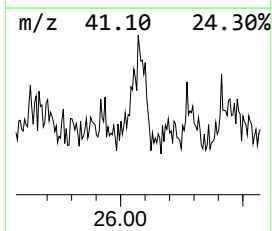
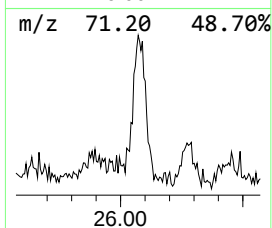
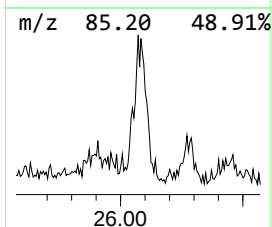
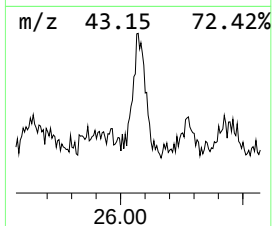
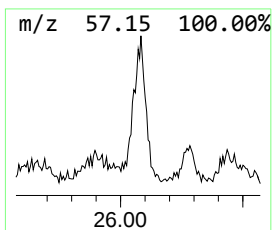
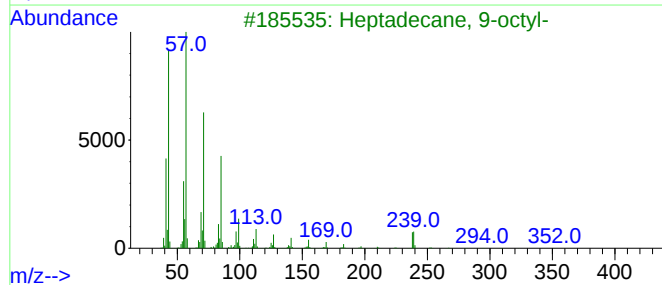
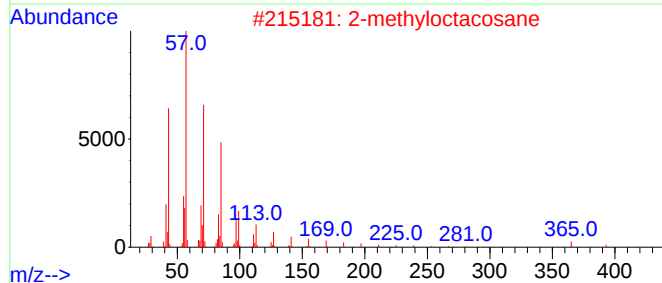
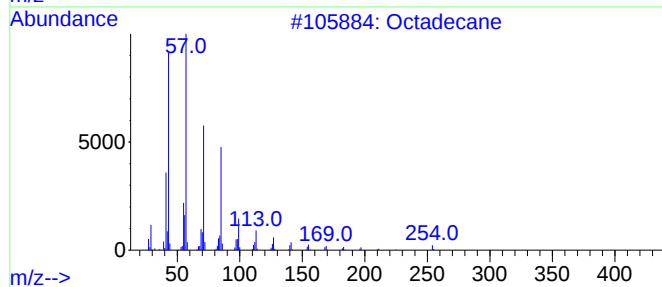
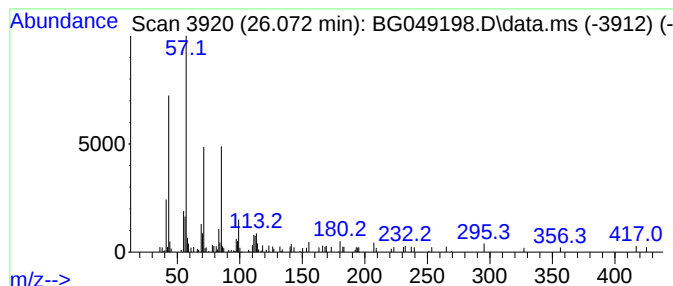
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG063021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Octadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.072	2.04 ng	70186	Perylene-d12	25.067

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	80
2			2-methyloctacosane	408	C29H60	1000376-72-8	80
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	80
4			Eicosane	282	C20H42	000112-95-8	80
5			Pentadecane	212	C15H32	000629-62-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070721\
 Data File : BG049198.D
 Acq On : 7 Jul 2021 14:00
 Operator : CG/JU
 Sample : M2933-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG063021.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
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unknown3.210	3.210	3.1	ng	36357	1	8.087	234167	20.0
n-Hexadecanoic ...	18.357	7.3	ng	190229	4	17.482	523539	20.0
Morpholine, 4-p...	18.704	9.8	ng	255816	4	17.482	523539	20.0
unknown18.839	18.839	16.4	ng	429785	4	17.482	523539	20.0
Diethylene glyc...	18.956	27.3	ng	714827	4	17.482	523539	20.0
Octadecanoic acid	19.620	10.7	ng	280346	4	17.482	523539	20.0
Octacosane	19.885	4.6	ng	137675	5	21.759	596186	20.0
Tetracosane	20.690	5.8	ng	173288	5	21.759	596186	20.0
Cyclotetradecane	21.401	8.0	ng	239102	5	21.759	596186	20.0
Eicosane	21.483	7.6	ng	225822	5	21.759	596186	20.0
Heptacosane	22.382	6.0	ng	179060	5	21.759	596186	20.0
Heneicosane	23.422	4.1	ng	141117	6	25.067	689322	20.0
Undecane	24.621	2.2	ng	75938	6	25.067	689322	20.0
Octadecane	26.072	2.0	ng	70186	6	25.067	689322	20.0