

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062521\
 Data File : BG049092.D
 Acq On : 25 Jun 2021 15:57
 Operator : CG/JU
 Sample : M2824-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ABN-1

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-BG061121.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG049092.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.157	15	20	29	rVB2	53071	120192	1.07%	0.301%
2	4.585	256	263	276	rVB	2152690	3500890	31.27%	8.775%
3	5.208	359	369	386	rVB	2277709	4308331	38.49%	10.799%
4	7.258	713	718	727	rVB	122323	218139	1.95%	0.547%
5	8.110	856	863	870	rVB2	146888	247049	2.21%	0.619%
6	9.279	1052	1062	1070	rBV	409550	755078	6.75%	1.893%
7	10.695	1296	1303	1313	rBV	66886	133437	1.19%	0.334%
8	10.930	1336	1343	1352	rVB	172030	315900	2.82%	0.792%
9	13.374	1752	1759	1766	rBV	953048	1475749	13.18%	3.699%
10	14.749	1986	1993	2003	rVB	258511	417692	3.73%	1.047%
11	17.499	2455	2461	2467	rBV	332089	520413	4.65%	1.304%
12	18.404	2593	2615	2629	rBV	853323	3754556	33.54%	9.411%
13	18.580	2629	2645	2655	rVV	1154367	4558645	40.72%	11.427%
14	18.739	2655	2672	2681	rVB	3444391	11194590	100.00%	28.060%
15	18.886	2688	2697	2710	rVB	272902	765978	6.84%	1.920%
16	19.644	2817	2826	2833	rVV3	95284	217215	1.94%	0.544%
17	19.732	2834	2841	2848	rVB2	188384	400623	3.58%	1.004%
18	20.108	2898	2905	2911	rVB	1774591	2392179	21.37%	5.996%
19	20.595	2981	2988	2995	rBV	313385	536944	4.80%	1.346%
20	21.412	3115	3127	3135	rBV	321756	694409	6.20%	1.741%
21	21.788	3184	3191	3196	rBV2	358051	610223	5.45%	1.530%
22	22.311	3273	3280	3291	rVB3	198847	520124	4.65%	1.304%
23	23.380	3450	3462	3471	rBV2	166637	462356	4.13%	1.159%
24	24.614	3663	3672	3681	rVB2	127874	381894	3.41%	0.957%
25	25.114	3745	3757	3767	rBV	202753	653316	5.84%	1.638%
26	26.077	3911	3921	3936	rVB4	76290	327673	2.93%	0.821%
27	27.822	4206	4218	4231	rVB3	55075	257358	2.30%	0.645%
28	28.950	4397	4410	4415	rBV3	35447	154048	1.38%	0.386%

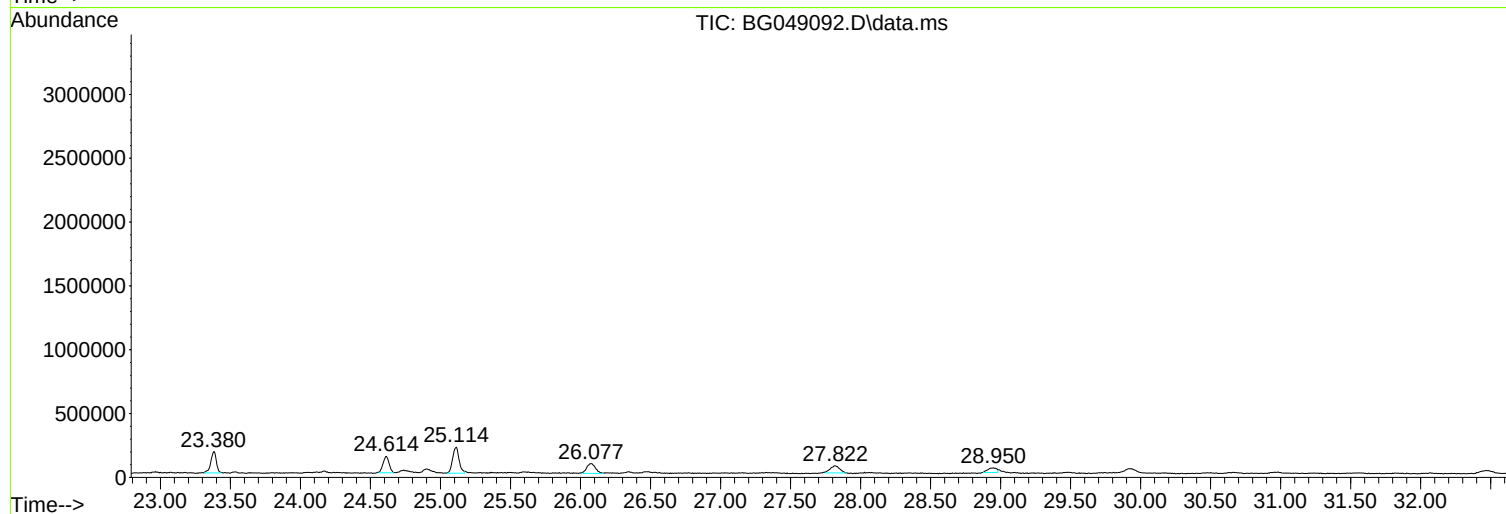
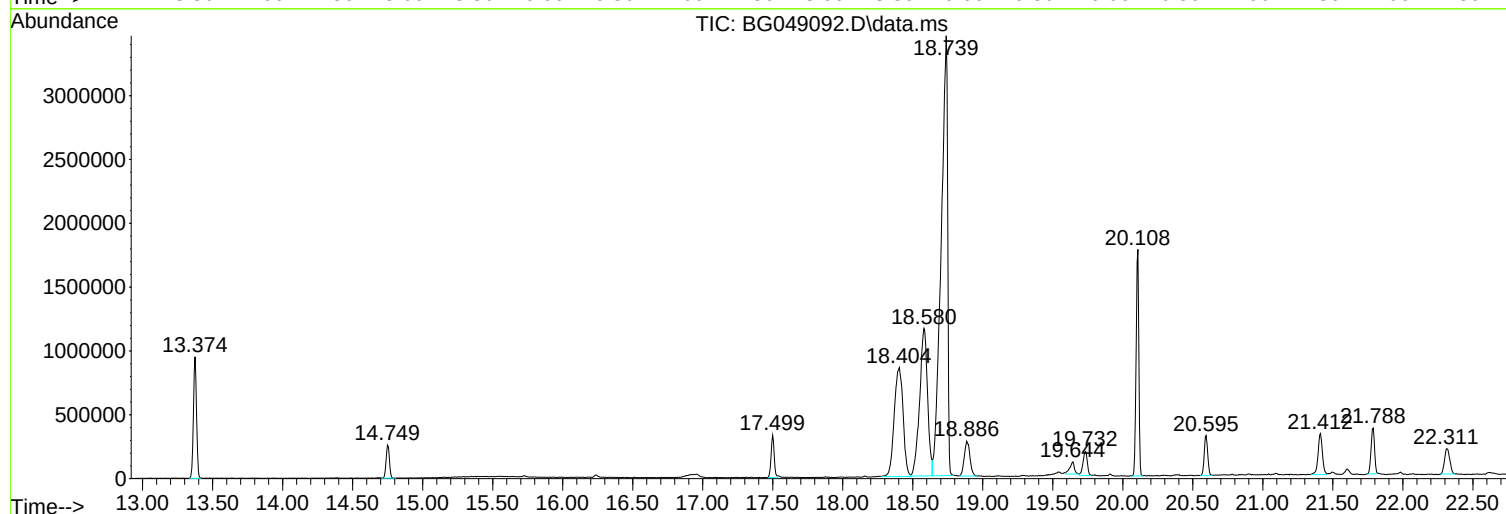
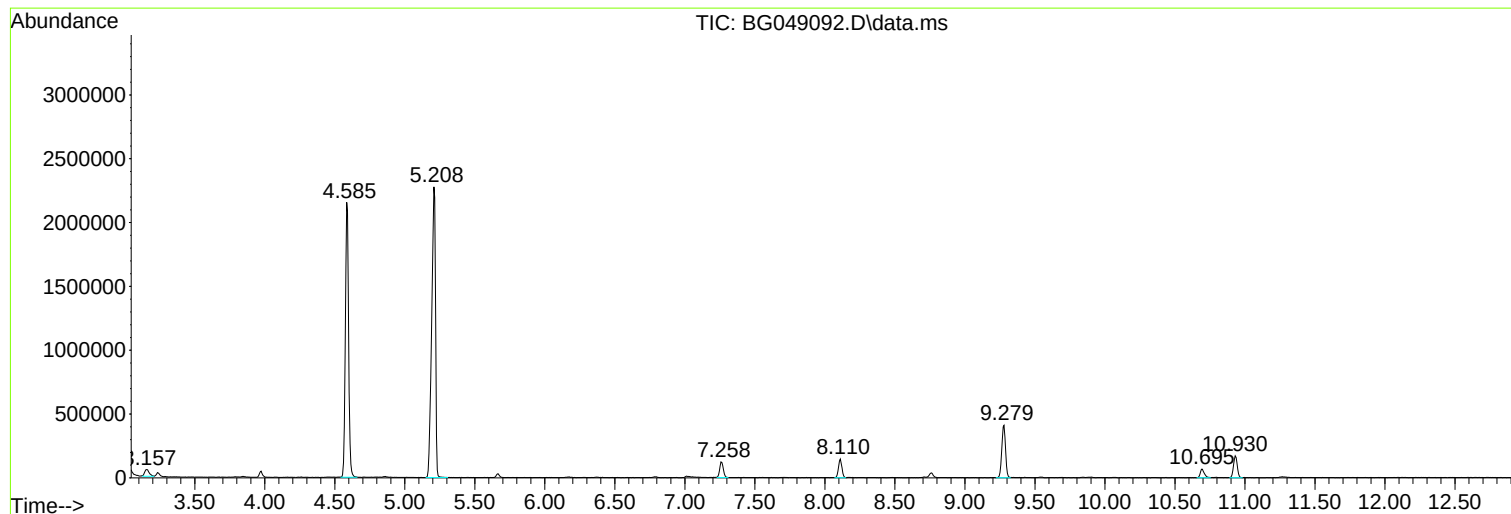
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Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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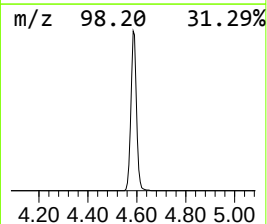
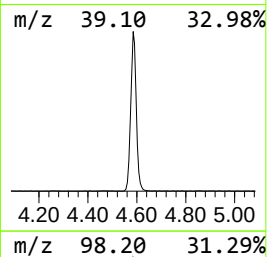
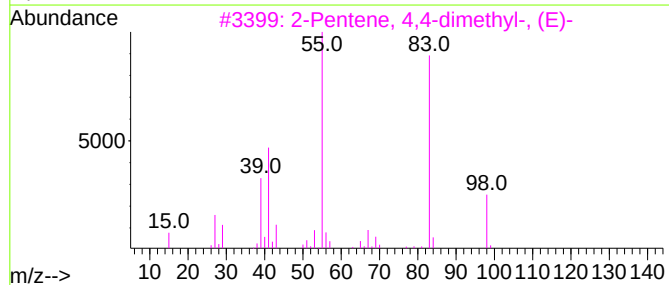
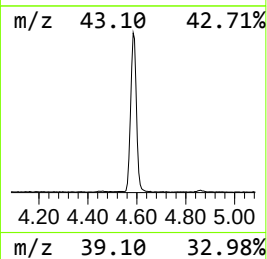
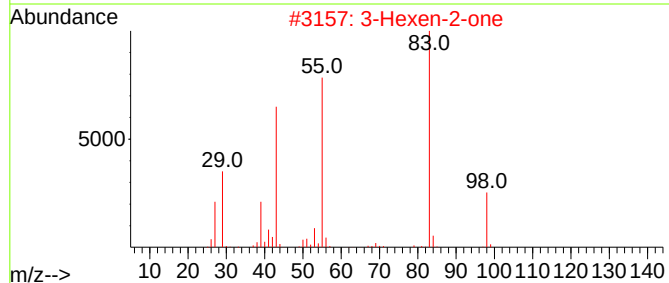
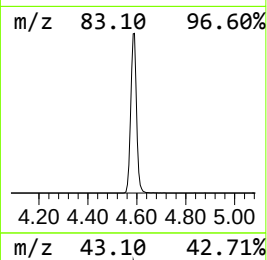
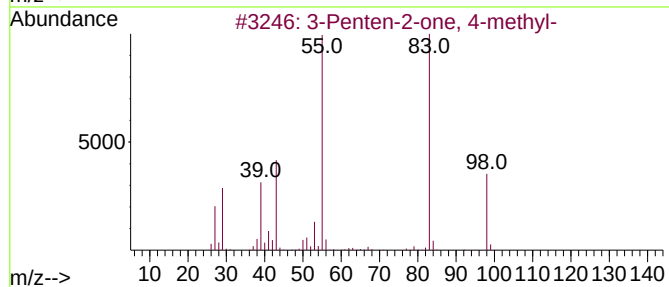
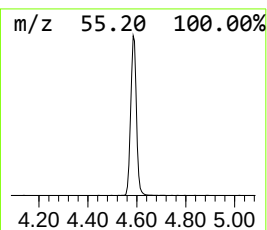
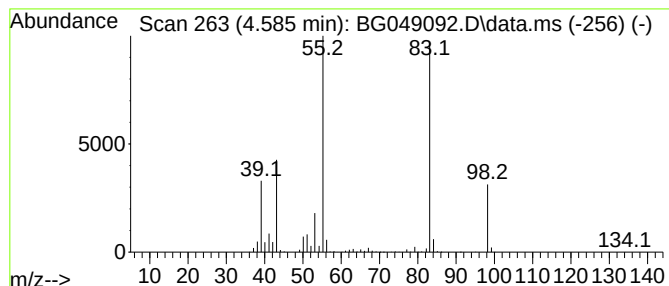
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.585	283.42 ng	3500890	1,4-Dichlorobenzene-d4	8.110

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2		3-Hexen-2-one	98	C6H10O	000763-93-9	91
3		2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	86
4		2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	86
5		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	80



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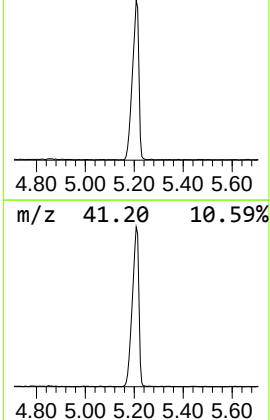
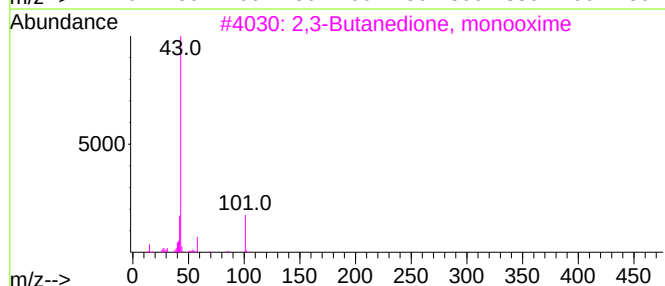
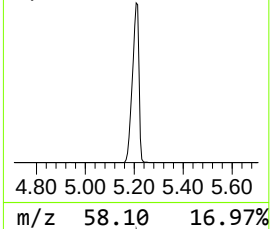
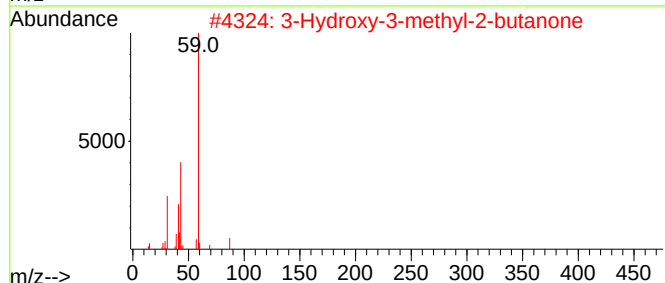
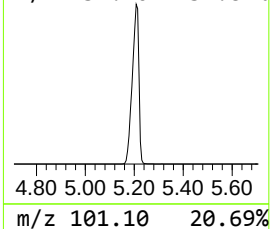
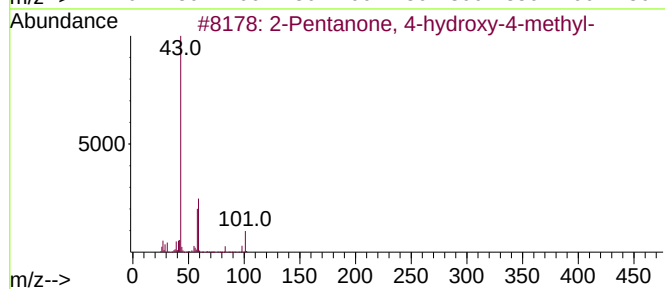
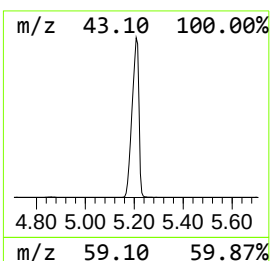
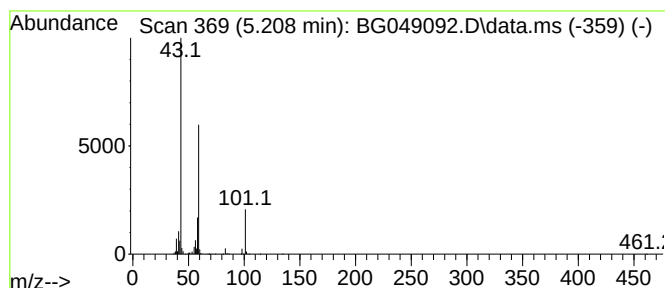
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.208	348.78 ng	4308330	1,4-Dichlorobenzene-d4	8.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	25
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
4			Butane	58	C4H10	000106-97-8	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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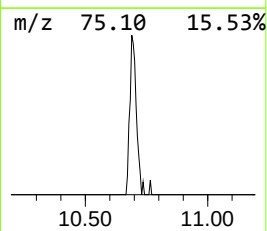
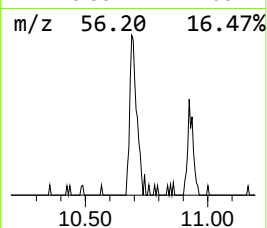
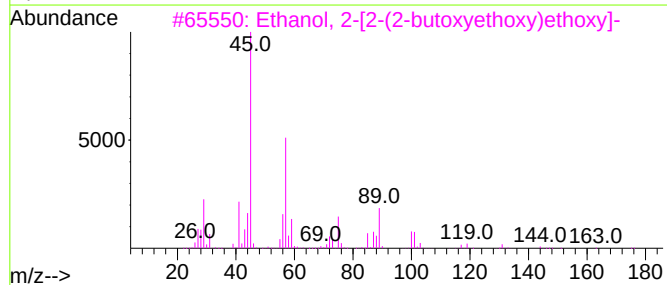
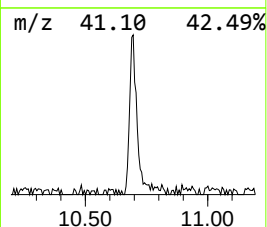
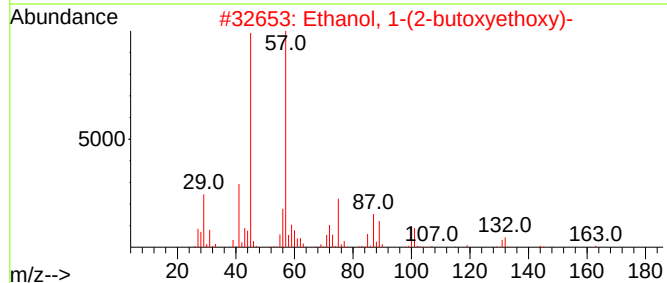
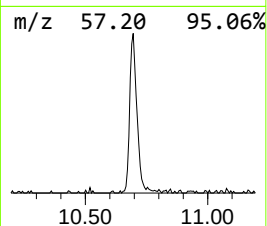
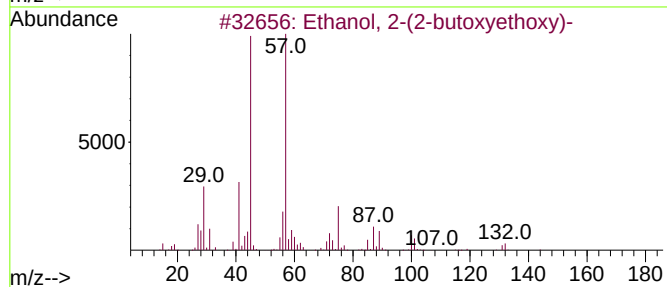
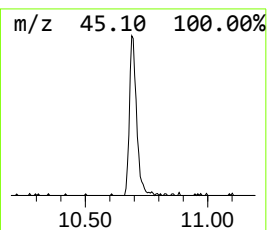
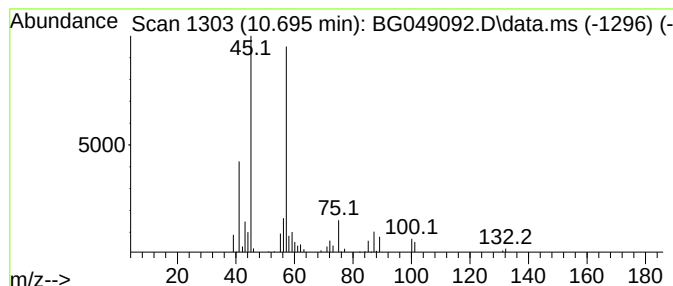
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.695	8.45 ng	133437	Naphthalene-d8	10.930

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3			Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	72
4			Methoxyacetic acid, heptyl ester	188	C10H20O3	168920-36-3	50
5			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	45



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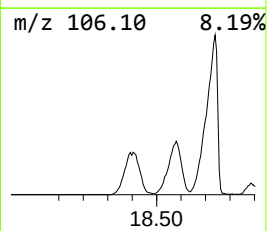
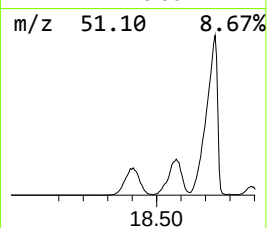
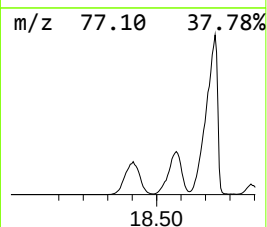
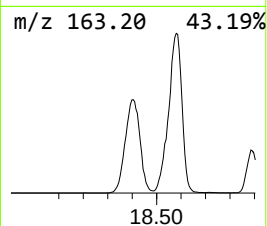
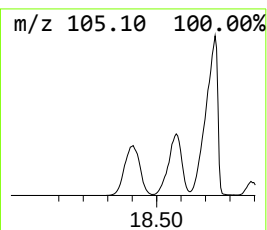
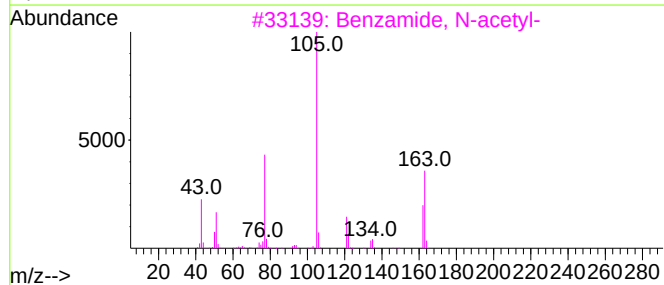
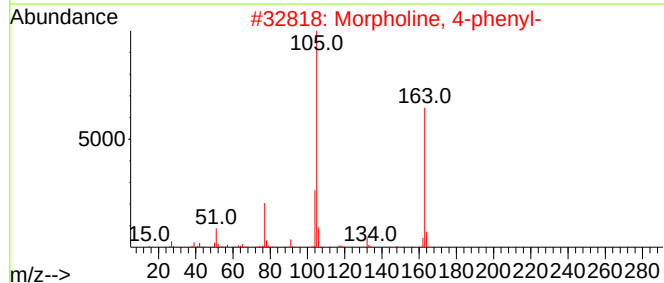
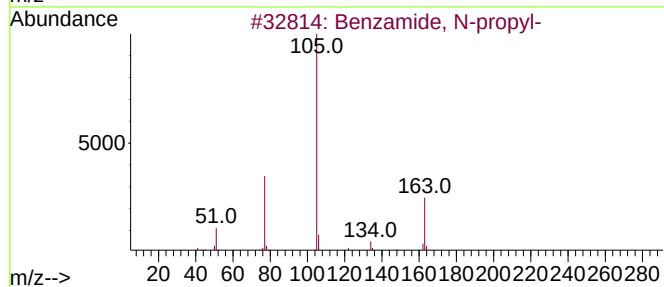
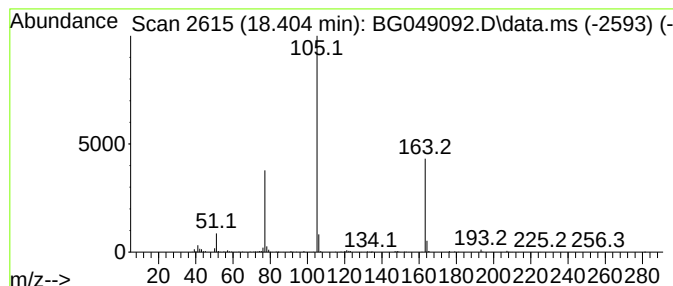
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzamide, N-propyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.404	144.29 ng	3754560	Phenanthrene-d10	17.499

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzamide, N-propyl-	163	C10H13NO	010546-70-0	72
2			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	64
3			Benzamide, N-acetyl-	163	C9H9NO2	001575-95-7	56
4			(3-Phenyloxiran-2-yl)methyl benz...	254	C16H14O3	1000366-96-1	42
5			Benzoic acid, pentafluorophenyl ...	288	C13H5F5O2	1000360-67-9	38



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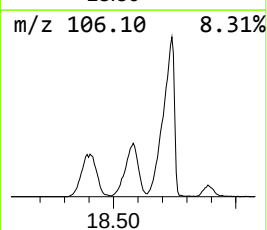
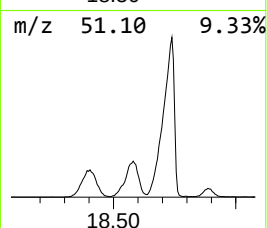
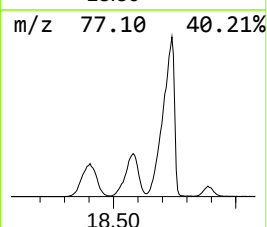
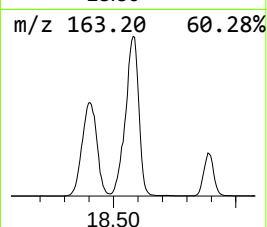
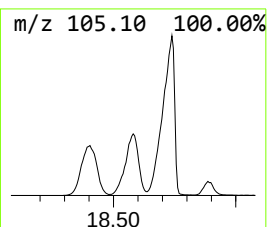
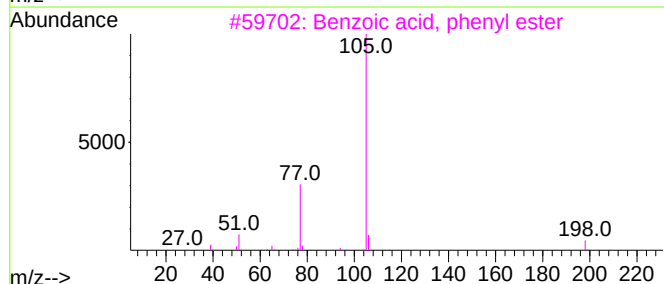
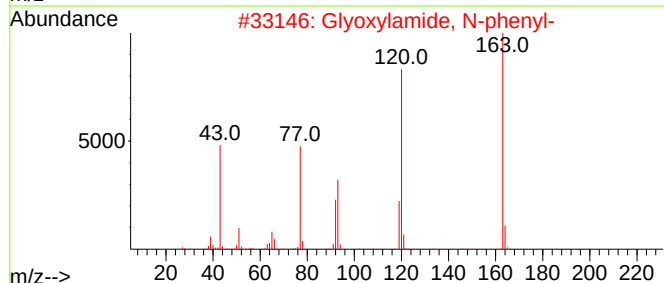
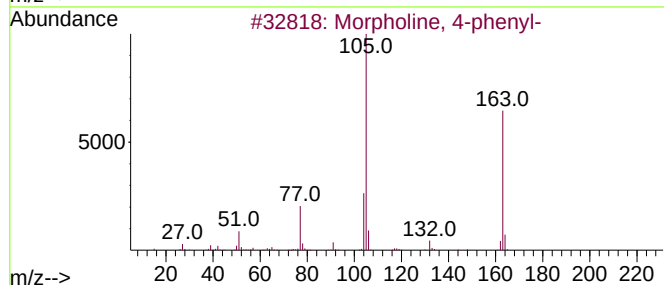
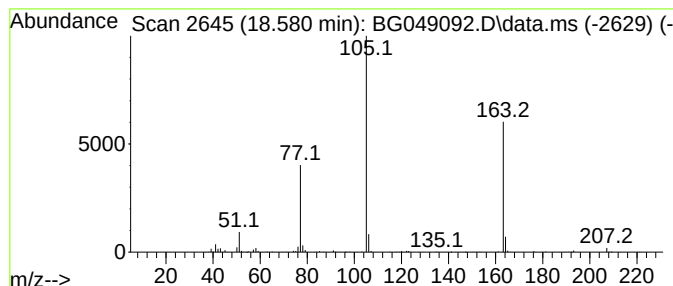
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Morpholine, 4-phenyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.580	175.19 ng	4558650	Phenanthrene-d10	17.499

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	53
2			Glyoxylamide, N-phenyl-	163	C9H9NO2	032331-52-5	53
3			Benzoic acid, phenyl ester	198	C13H10O2	000093-99-2	47
4			Benzoic acid, 3,5-difluophenyl e...	234	C13H8F2O2	1000357-79-5	47
5			Benzeneacetic acid, .alpha.-oxo-...	178	C10H10O3	001603-79-8	47



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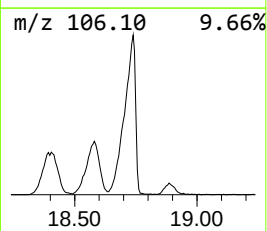
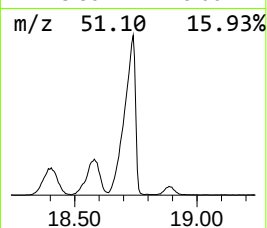
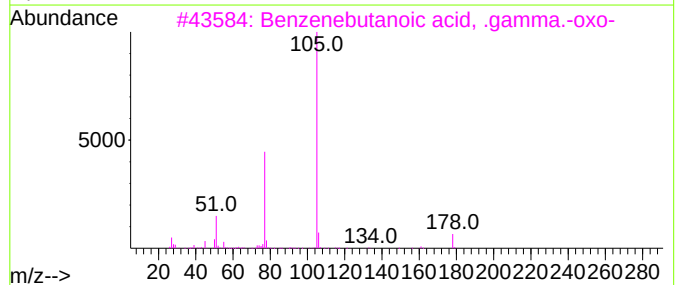
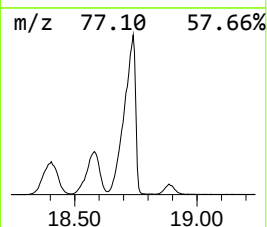
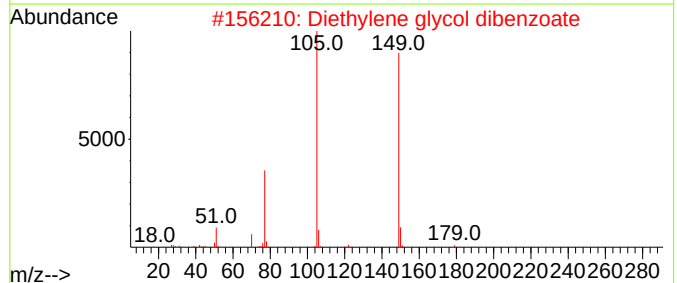
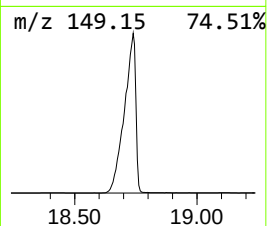
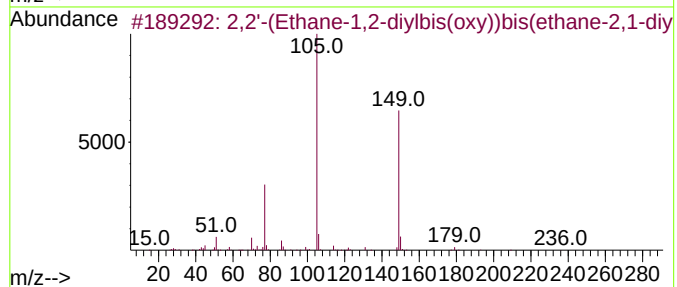
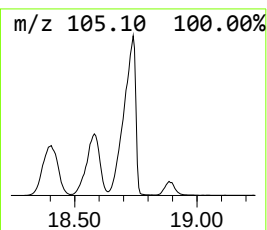
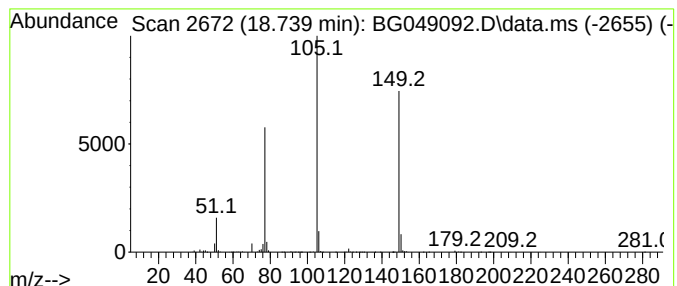
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 2,2'-(Ethane-1,2-diylbis(ox... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.739	430.22 ng	11194600	Phenanthrene-d10	17.499

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2'-(Ethane-1,2-diylbis(oxy))bi...	358	C20H22O6	1000366-99-4	74		
2	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	72		
3	Benzenebutanoic acid, .gamma.-oxo-	178	C10H10O3	002051-95-8	38		
4	2-Phenyl-4-ethylidene-2-oxazolin...	187	C11H9NO2	037791-41-6	38		
5	Benzoyl bromide	184	C7H5BrO	000618-32-6	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062521\
Data File : BG049092.D
Acq On : 25 Jun 2021 15:57
Operator : CG/JU
Sample : M2824-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

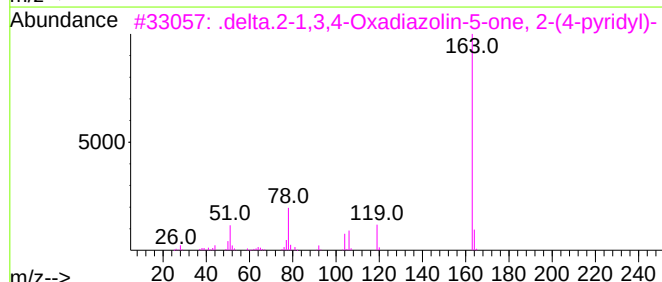
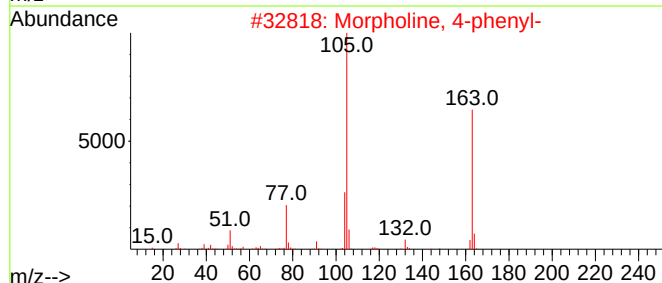
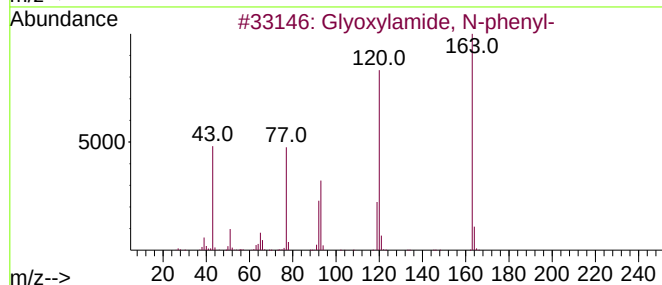
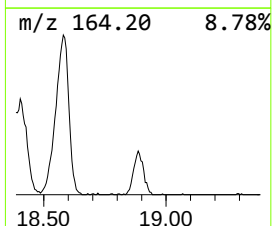
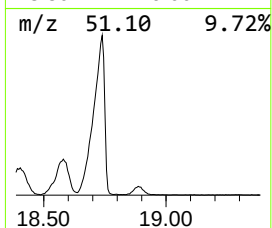
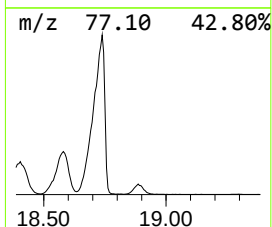
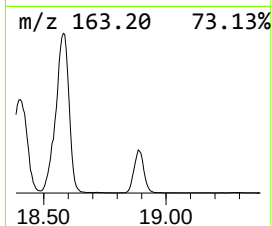
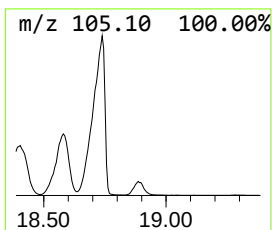
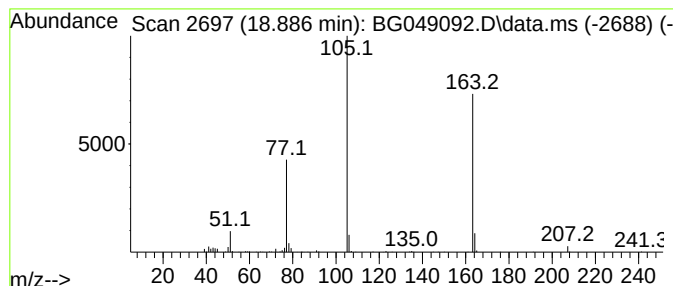
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Glyoxylamide, N-phenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD			R.T.
18.886	29.44 ng	765978	Phenanthrene-d10			17.499
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Glyoxylamide, N-phenyl-		163	C9H9NO2	032331-52-5	64
2	Morpholine, 4-phenyl-		163	C10H13NO	000092-53-5	59
3	.delta.2-1,3,4-Oxadiazolin-5-one...		163	C7H5N3O2	002845-82-1	47
4	Benzoyl isothiocyanate		163	C8H5NOS	000532-55-8	35
5	1,2-Bis(benzoyloxy)benzene		318	C20H14O4	000643-94-7	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062521\
Data File : BG049092.D
Acq On : 25 Jun 2021 15:57
Operator : CG/JU
Sample : M2824-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

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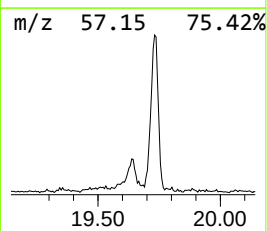
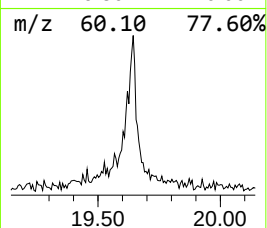
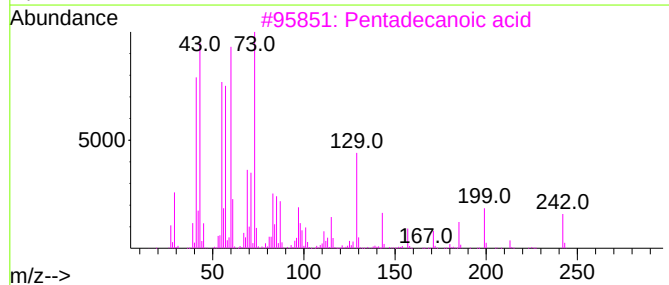
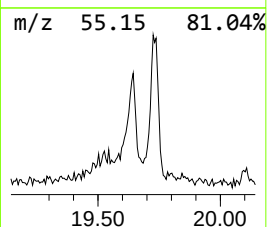
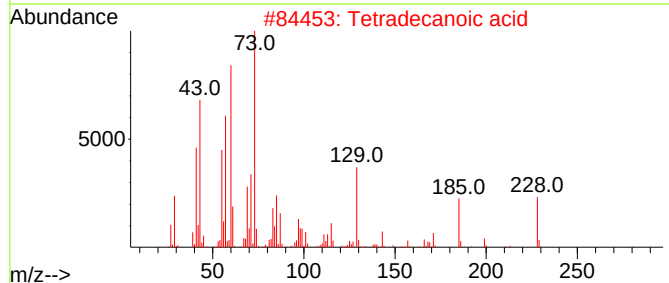
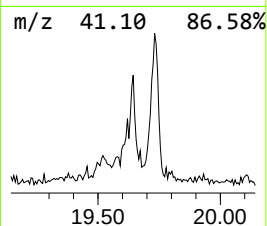
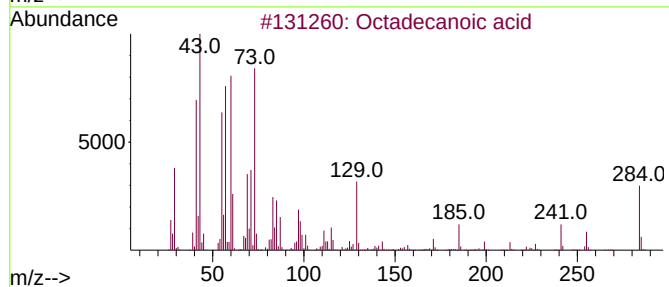
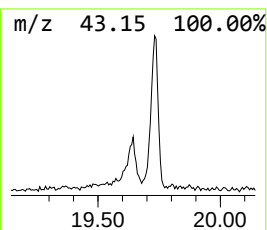
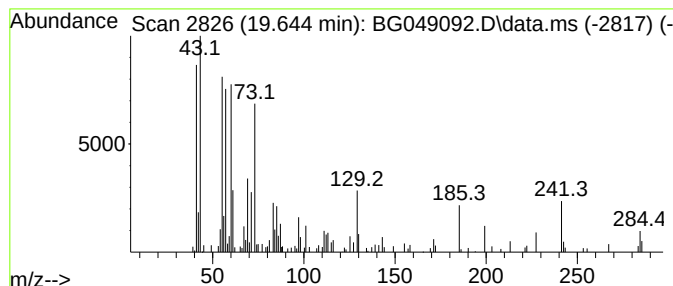
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.644	7.12 ng	217215	Chrysene-d12	21.788

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	62
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	53
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	52
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	49
5			Dodecanoic acid	200	C12H24O2	000143-07-7	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062521\
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Acq On : 25 Jun 2021 15:57
Operator : CG/JU
Sample : M2824-02
Misc :
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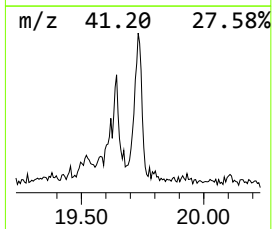
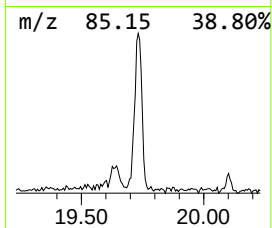
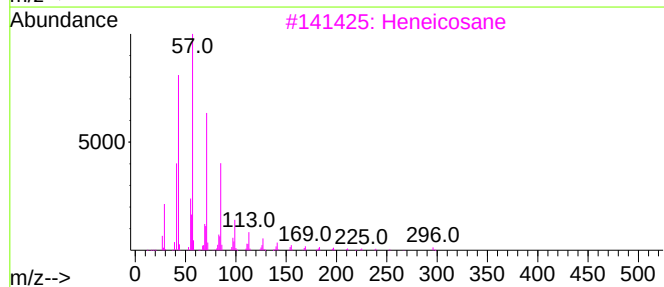
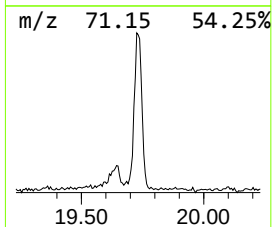
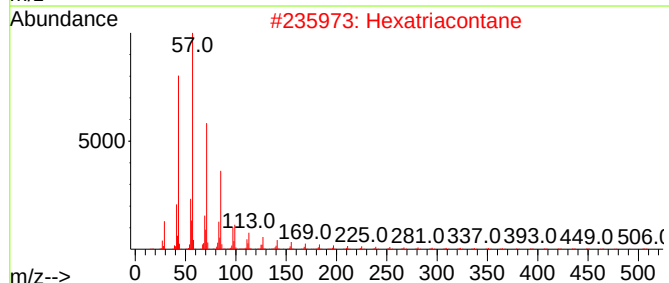
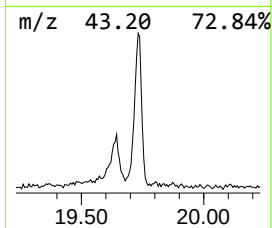
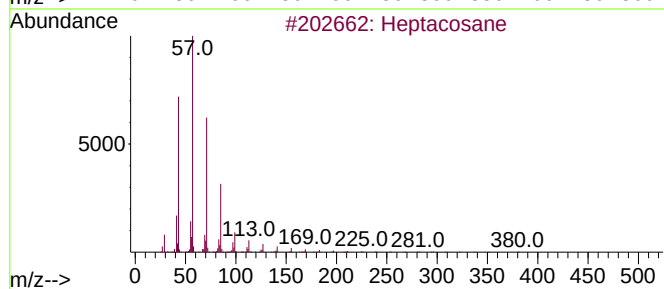
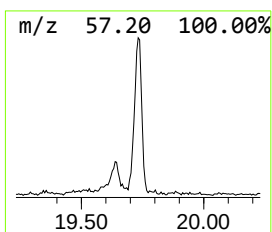
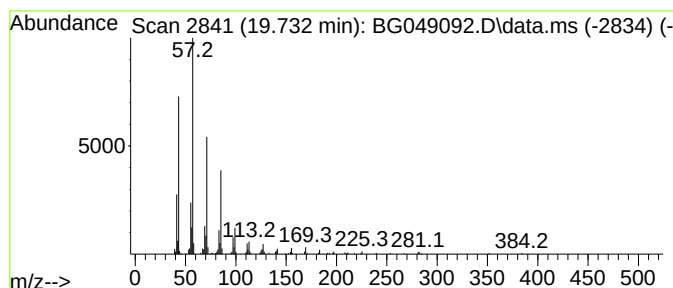
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.732	13.13 ng	400623	Chrysene-d12	21.788	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	91
2	Hexatriacontane	507	C36H74	000630-06-8	91
3	Heneicosane	296	C21H44	000629-94-7	90
4	Hexacosane	366	C26H54	000630-01-3	90
5	Eicosane	282	C20H42	000112-95-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062521\
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Misc :
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Instrument :
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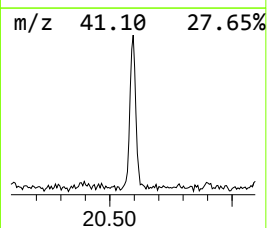
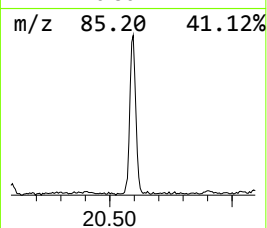
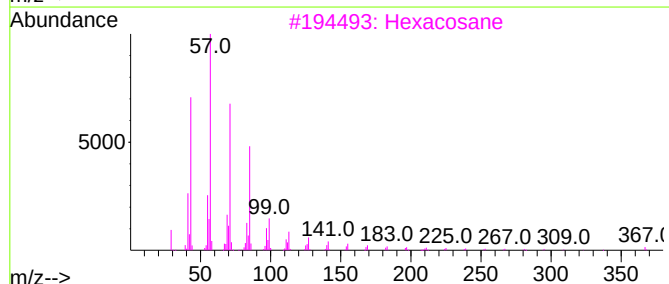
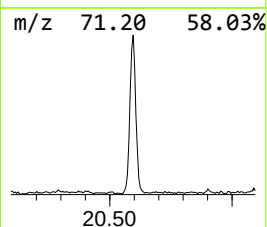
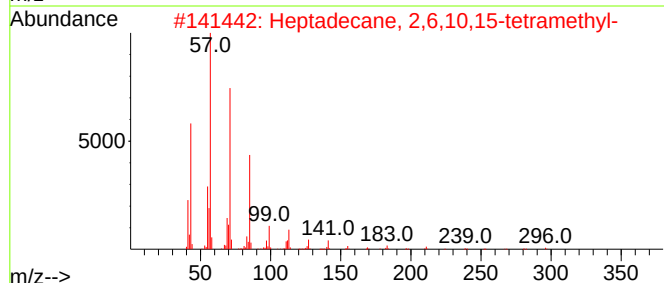
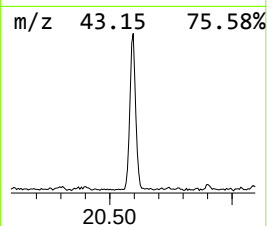
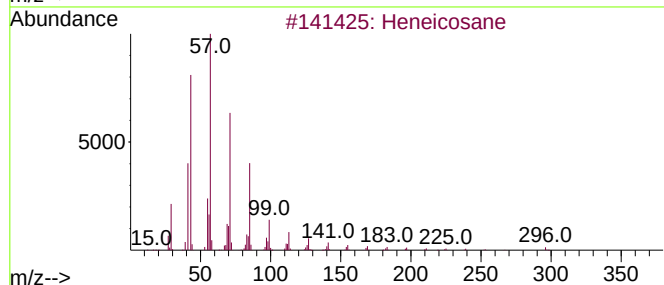
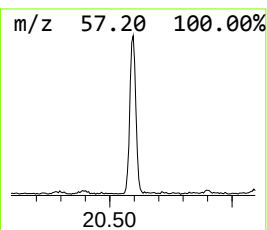
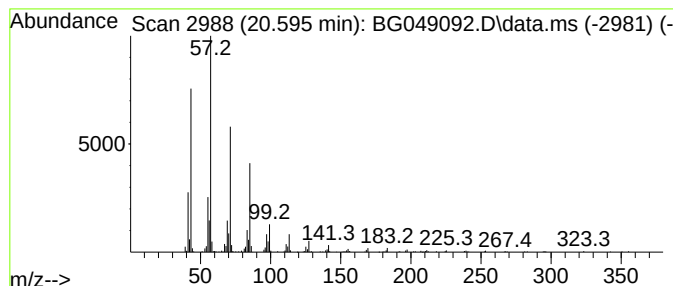
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.595	17.60 ng	536944	Chrysene-d12	21.788

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	96
2			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	96
3			Hexacosane	366	C26H54	000630-01-3	94
4			Hentriacontane	437	C31H64	000630-04-6	91
5			Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062521\
Data File : BG049092.D
Acq On : 25 Jun 2021 15:57
Operator : CG/JU
Sample : M2824-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

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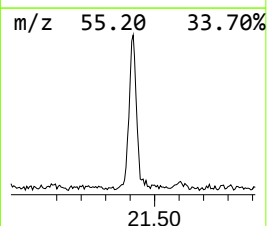
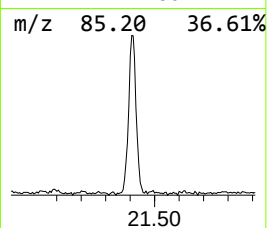
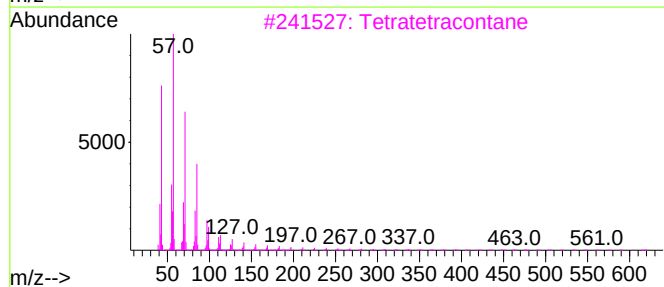
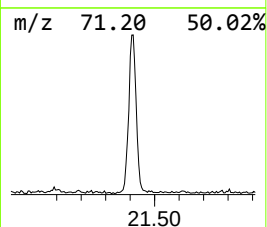
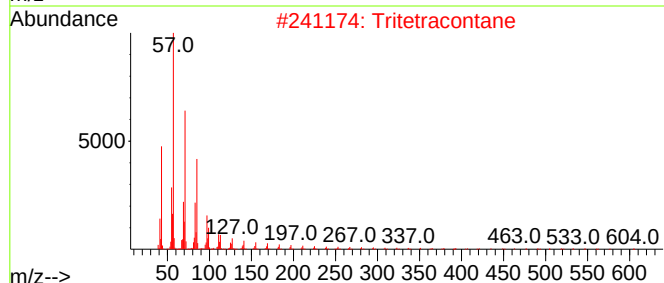
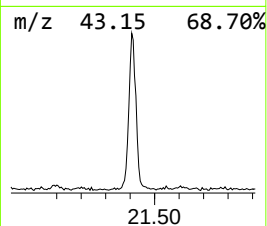
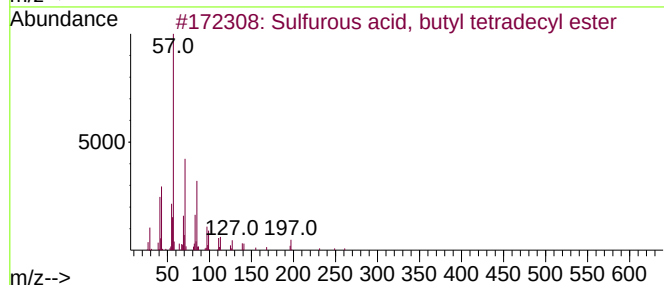
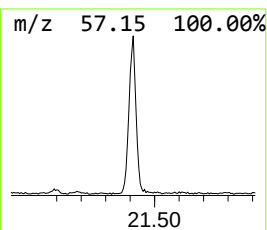
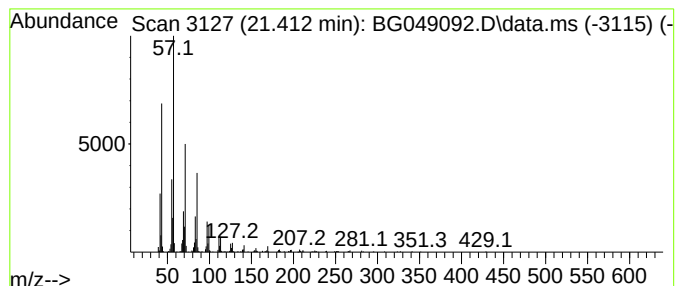
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Sulfurous acid, butyl tetra... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.412	22.76 ng	694409	Chrysene-d12	21.788

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1000309-18-1	91
2			Tritetracontane	605	C43H88	007098-21-7	90
3			Tetratetracontane	619	C44H90	007098-22-8	90
4			Heneicosane	296	C21H44	000629-94-7	87
5			Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062521\
Data File : BG049092.D
Acq On : 25 Jun 2021 15:57
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Misc :
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Instrument :
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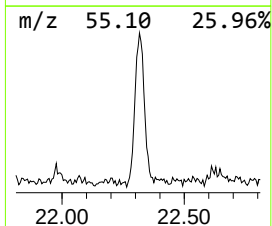
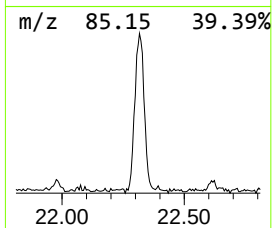
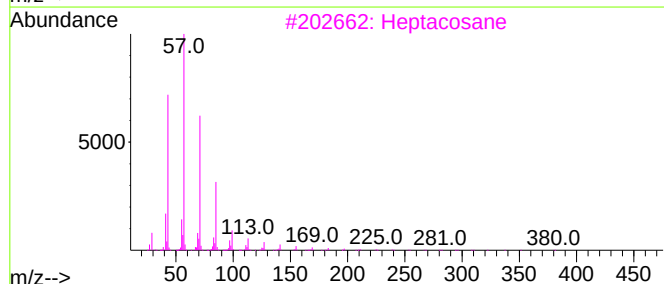
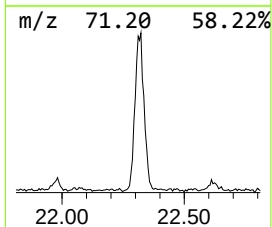
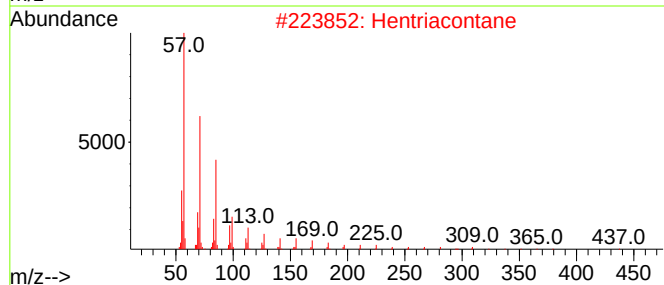
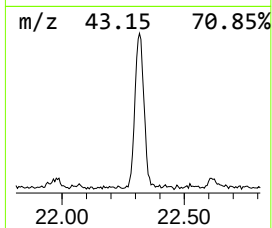
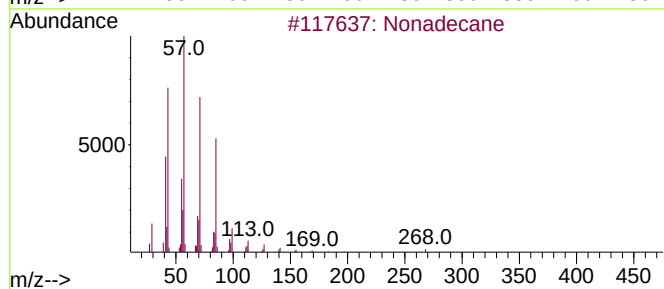
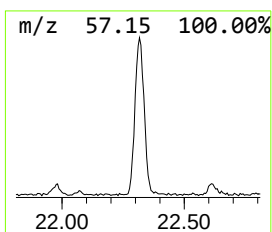
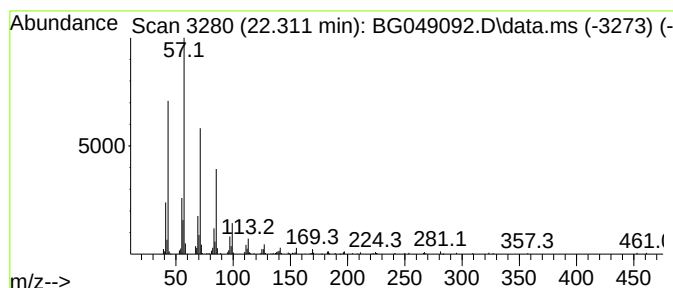
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Nonadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.311	17.05 ng	520124	Chrysene-d12	21.788

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	93
2		Hentriacontane	437	C31H64	000630-04-6	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062521\
Data File : BG049092.D
Acq On : 25 Jun 2021 15:57
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Sample : M2824-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
ABN-1

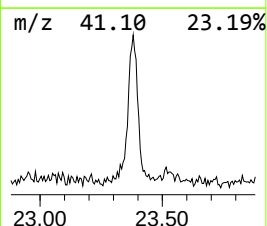
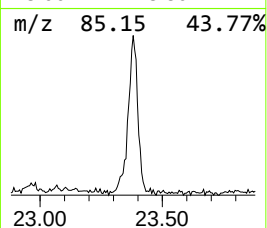
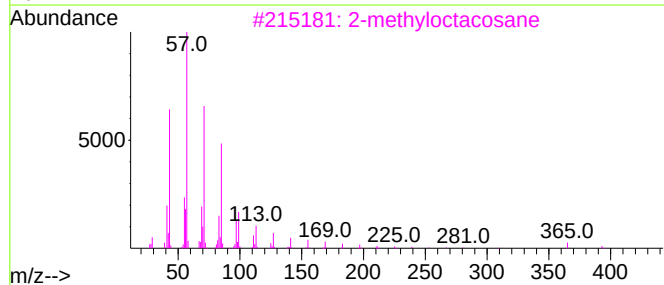
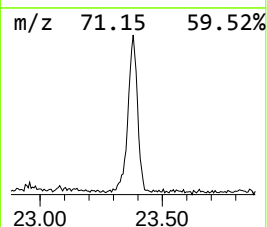
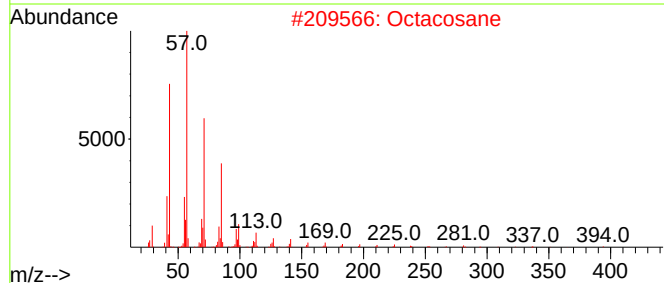
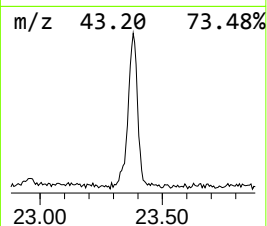
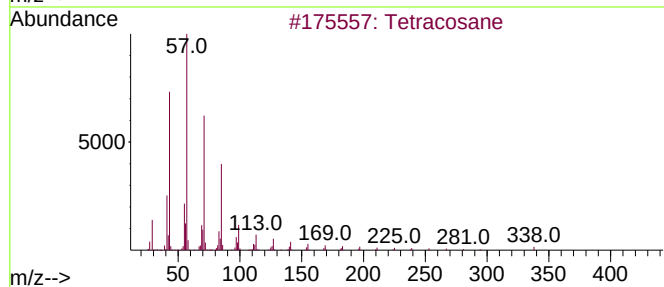
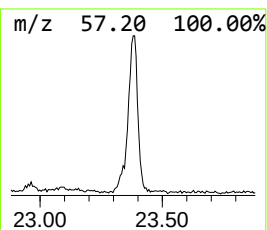
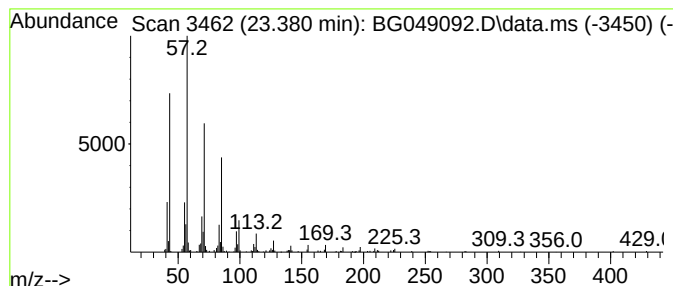
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Tetracosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.380	15.15 ng	462356	Chrysene-d12	21.788

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	90
2			Octacosane	394	C28H58	000630-02-4	90
3			2-methyloctacosane	408	C29H60	1000376-72-8	87
4			Heneicosane	296	C21H44	000629-94-7	83
5			Hexacosane	366	C26H54	000630-01-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062521\
Data File : BG049092.D
Acq On : 25 Jun 2021 15:57
Operator : CG/JU
Sample : M2824-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ABN-1

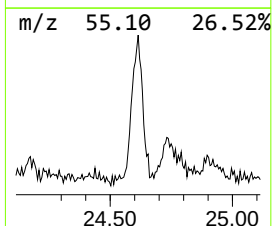
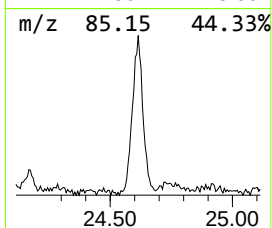
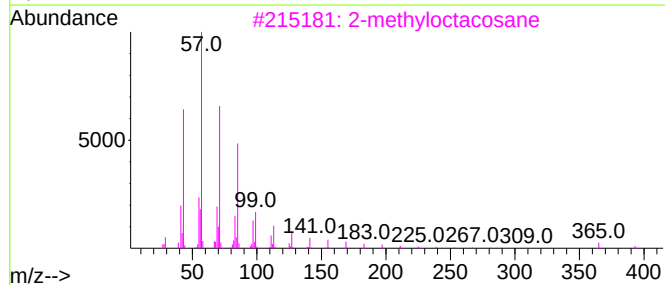
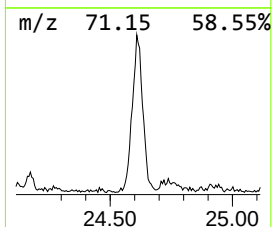
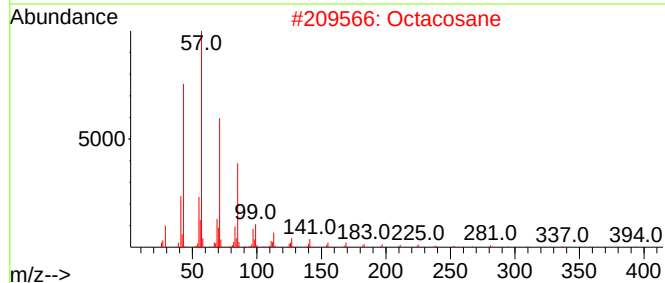
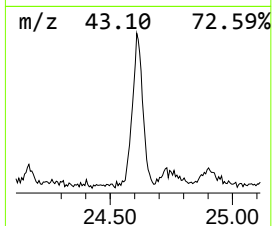
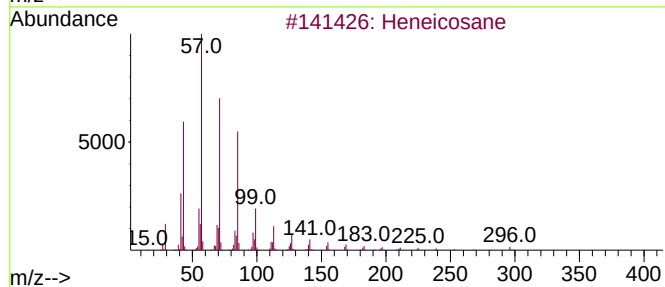
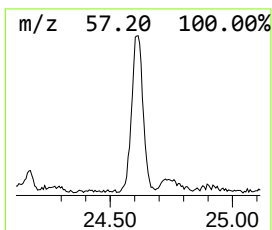
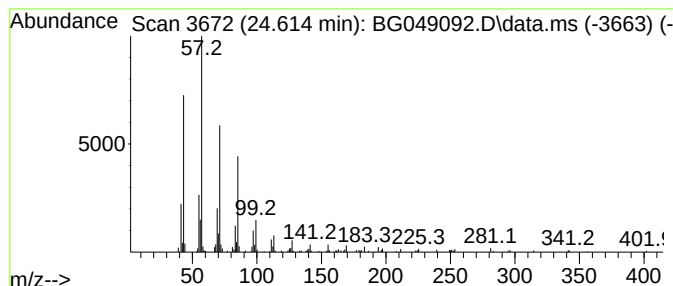
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Octacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.614	11.69 ng	381894	Perylene-d12	25.114

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	97
2			Octacosane	394	C28H58	000630-02-4	90
3			2-methyloctacosane	408	C29H60	1000376-72-8	90
4			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	90
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062521\
Data File : BG049092.D
Acq On : 25 Jun 2021 15:57
Operator : CG/JU
Sample : M2824-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

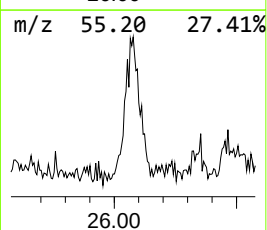
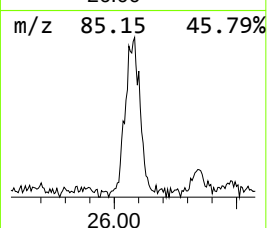
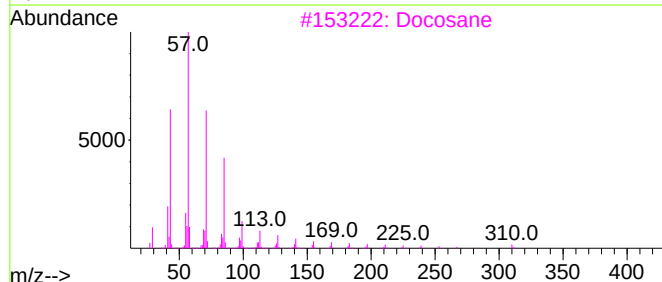
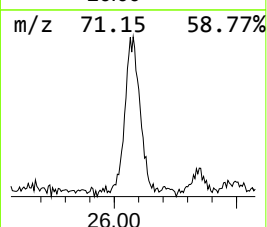
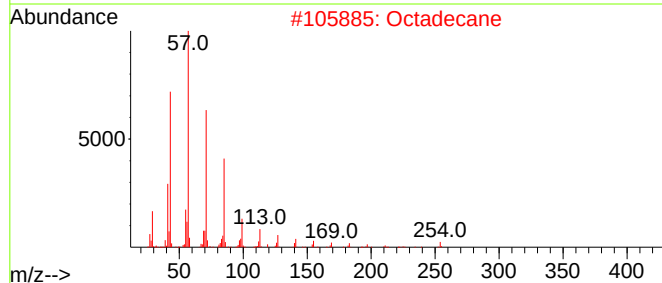
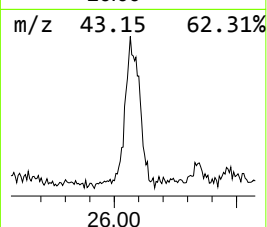
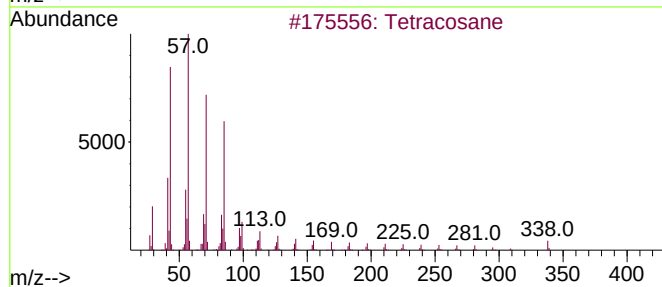
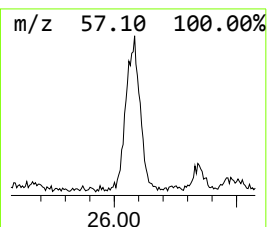
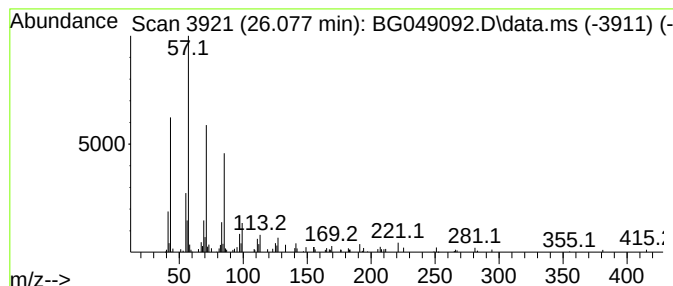
Instrument :
BNA_G
ClientSampleId :
ABN-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Octadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
26.077	10.03 ng	327673	Perylene-d12		25.114	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetracosane		338	C24H50	000646-31-1	87
2	Octadecane		254	C18H38	000593-45-3	83
3	Docosane		310	C22H46	000629-97-0	83
4	Eicosane		282	C20H42	000112-95-8	80
5	2-methyloctacosane		408	C29H60	1000376-72-8	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062521\
Data File : BG049092.D
Acq On : 25 Jun 2021 15:57
Operator : CG/JU
Sample : M2824-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
ABN-1

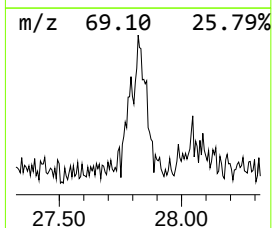
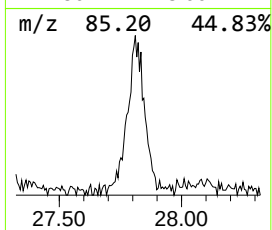
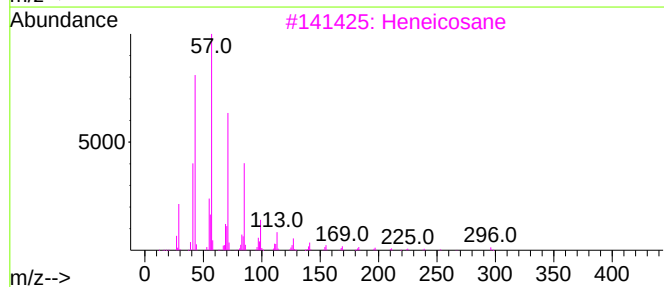
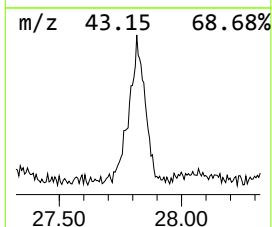
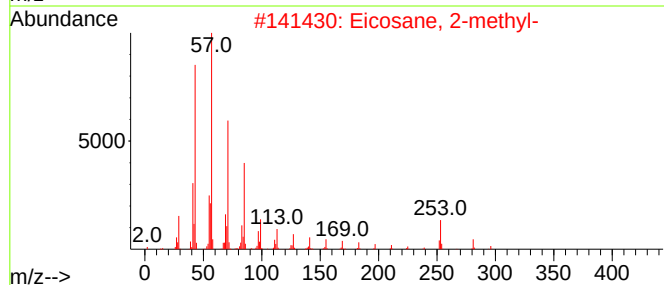
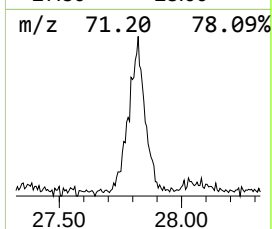
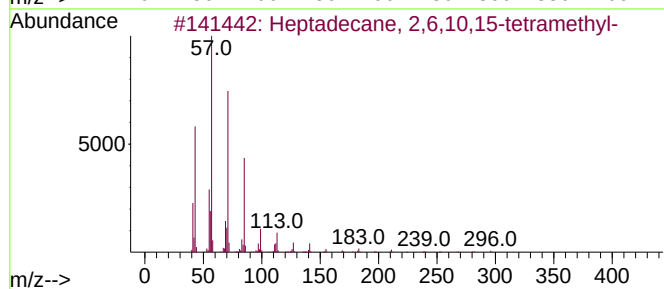
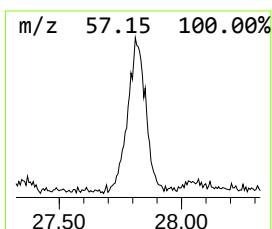
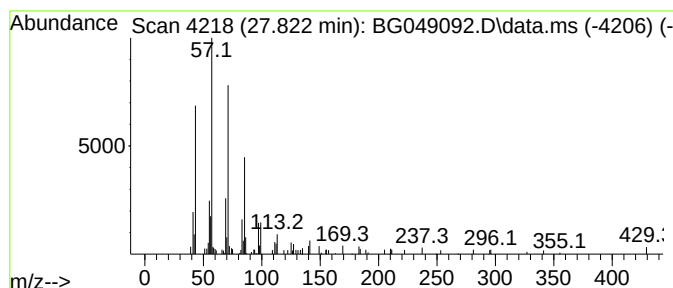
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Heptadecane, 2,6,10,15-tetr... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.822	7.88 ng	257358	Perylene-d12	25.114

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
2			Eicosane, 2-methyl-	296	C21H44	001560-84-5	93
3			Heneicosane	296	C21H44	000629-94-7	93
4			Docosane, 11-butyl-	366	C26H54	013475-76-8	91
5			Docosane	310	C22H46	000629-97-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062521\
Data File : BG049092.D
Acq On : 25 Jun 2021 15:57
Operator : CG/JU
Sample : M2824-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

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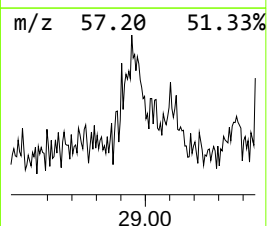
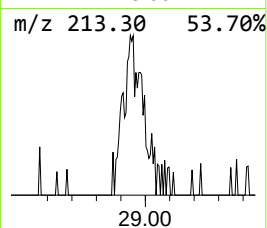
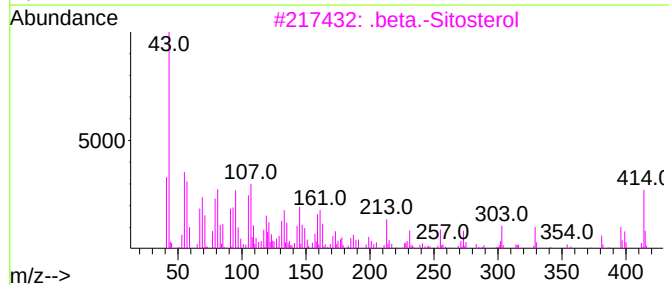
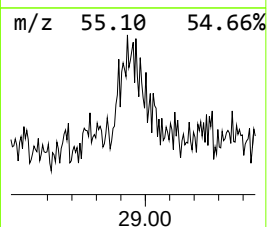
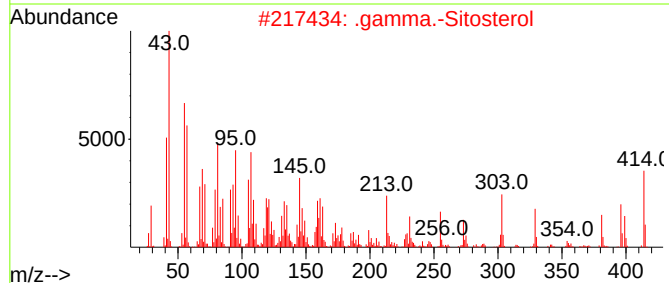
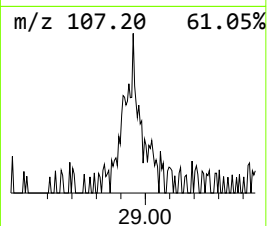
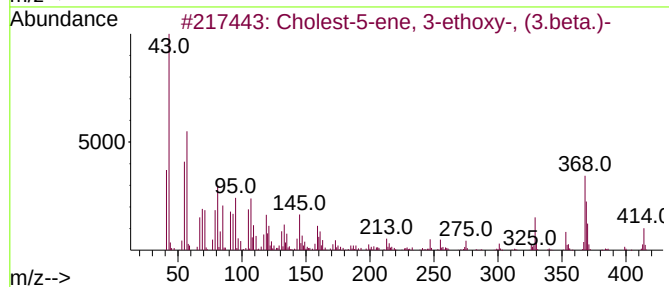
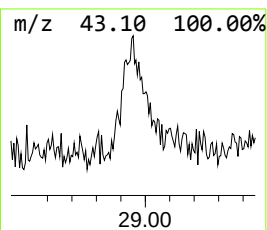
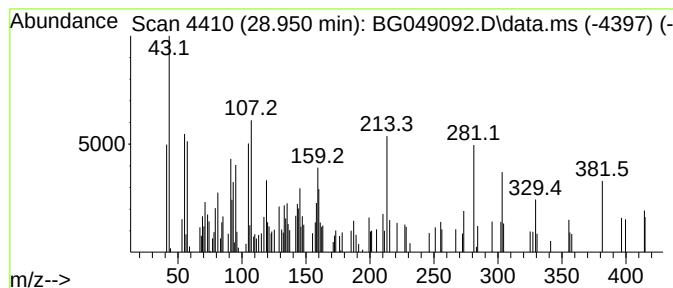
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 unknown28.950 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.950	4.72 ng	154048	Perylene-d12	25.114

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cholest-5-ene, 3-ethoxy-, (3.bet...	414	C29H50O	000986-19-6	25
2		.gamma.-Sitosterol	414	C29H50O	000083-47-6	22
3		.beta.-Sitosterol	414	C29H50O	000083-46-5	14
4		1-Oxaspiro[4.4]nonane-4-carboxam...	273	C16H19NO3	1000319-97-0	11
5		Ethinyl Estradiol	296	C20H24O2	000057-63-6	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062521\
 Data File : BG049092.D
 Acq On : 25 Jun 2021 15:57
 Operator : CG/JU
 Sample : M2824-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ABN-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.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
3-Penten-2-one,...	4.585	283.4	ng	3500890	1	8.110	247049	20.0
2-Pentanone, 4-...	5.208	348.8	ng	4308330	1	8.110	247049	20.0
Ethanol, 2-(2-b...	10.695	8.4	ng	133437	2	10.930	315900	20.0
Benzamide, N-pr...	18.404	144.3	ng	3754560	4	17.499	520413	20.0
Morpholine, 4-p...	18.580	175.2	ng	4558650	4	17.499	520413	20.0
2,2'-(Ethane-1,...	18.739	430.2	ng	11194600	4	17.499	520413	20.0
Glyoxylamide, N...	18.886	29.4	ng	765978	4	17.499	520413	20.0
Octadecanoic acid	19.644	7.1	ng	217215	5	21.788	610223	20.0
Heptacosane	19.732	13.1	ng	400623	5	21.788	610223	20.0
Heneicosane	20.595	17.6	ng	536944	5	21.788	610223	20.0
Sulfurous acid,...	21.412	22.8	ng	694409	5	21.788	610223	20.0
Nonadecane	22.311	17.1	ng	520124	5	21.788	610223	20.0
Tetracosane	23.380	15.2	ng	462356	5	21.788	610223	20.0
Octacosane	24.614	11.7	ng	381894	6	25.114	653316	20.0
Octadecane	26.077	10.0	ng	327673	6	25.114	653316	20.0
Heptadecane, 2,...	27.822	7.9	ng	257358	6	25.114	653316	20.0
unknown28.950	28.950	4.7	ng	154048	6	25.114	653316	20.0