

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
 Data File : BM029025.D
 Acq On : 08 Mar 2021 18:55
 Operator : CG/JU
 Sample : M1575-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 IDW-GW806-AQ-030321

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_M\Methods\8270-BM022321.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM029025.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.363	61	64	70	rBV	13231	17667	2.99%	0.662%
2	4.775	300	304	312	rVB	5036	6825	1.16%	0.256%
3	5.246	379	384	392	rBB	89877	126613	21.44%	4.746%
4	5.722	460	465	470	rBV	18824	25926	4.39%	0.972%
5	6.722	631	635	643	rBV2	3908	7798	1.32%	0.292%
6	6.869	655	660	670	rBB	53418	87486	14.81%	3.279%
7	7.669	791	796	802	rBB	37335	55302	9.36%	2.073%
8	7.740	803	808	815	rBB	6696	9795	1.66%	0.367%
9	8.310	899	905	910	rBB	5113	8723	1.48%	0.327%
10	8.845	989	996	1005	rVB	108071	169336	28.67%	6.347%
11	9.828	1158	1163	1167	rVB3	6273	8834	1.50%	0.331%
12	9.886	1168	1173	1183	rBB	9693	14456	2.45%	0.542%
13	10.463	1265	1271	1277	rBV	47483	73384	12.42%	2.751%
14	11.363	1418	1424	1432	rVB2	7880	13270	2.25%	0.497%
15	12.951	1688	1694	1701	rVB	266513	365857	61.94%	13.714%
16	13.804	1835	1839	1844	rVB	9845	12210	2.07%	0.458%
17	14.221	1903	1910	1915	rBV3	3551	7324	1.24%	0.275%
18	14.292	1917	1922	1925	rVV	4327	6352	1.08%	0.238%
19	14.339	1925	1930	1936	rVB2	64339	89639	15.18%	3.360%
20	15.833	2178	2184	2192	rBV	210963	294781	49.91%	11.050%
21	17.086	2391	2397	2402	rBV	77545	103022	17.44%	3.862%
22	17.862	2519	2529	2534	rBV4	3972	7900	1.34%	0.296%
23	17.980	2544	2549	2560	rBV	41108	54491	9.23%	2.043%
24	18.721	2670	2675	2681	rVB	22627	26825	4.54%	1.006%
25	18.809	2686	2690	2694	rVB2	4965	5964	1.01%	0.224%
26	19.262	2763	2767	2772	rBV	23887	28873	4.89%	1.082%
27	19.709	2838	2843	2848	rBV	500034	590664	100.00%	22.141%
28	20.221	2924	2930	2943	rBV	89267	130246	22.05%	4.882%
29	20.974	3053	3058	3066	rBV	53419	67676	11.46%	2.537%
30	21.256	3102	3106	3111	rBV	104498	125715	21.28%	4.712%
31	23.415	3468	3473	3481	rVB2	72239	124843	21.14%	4.680%

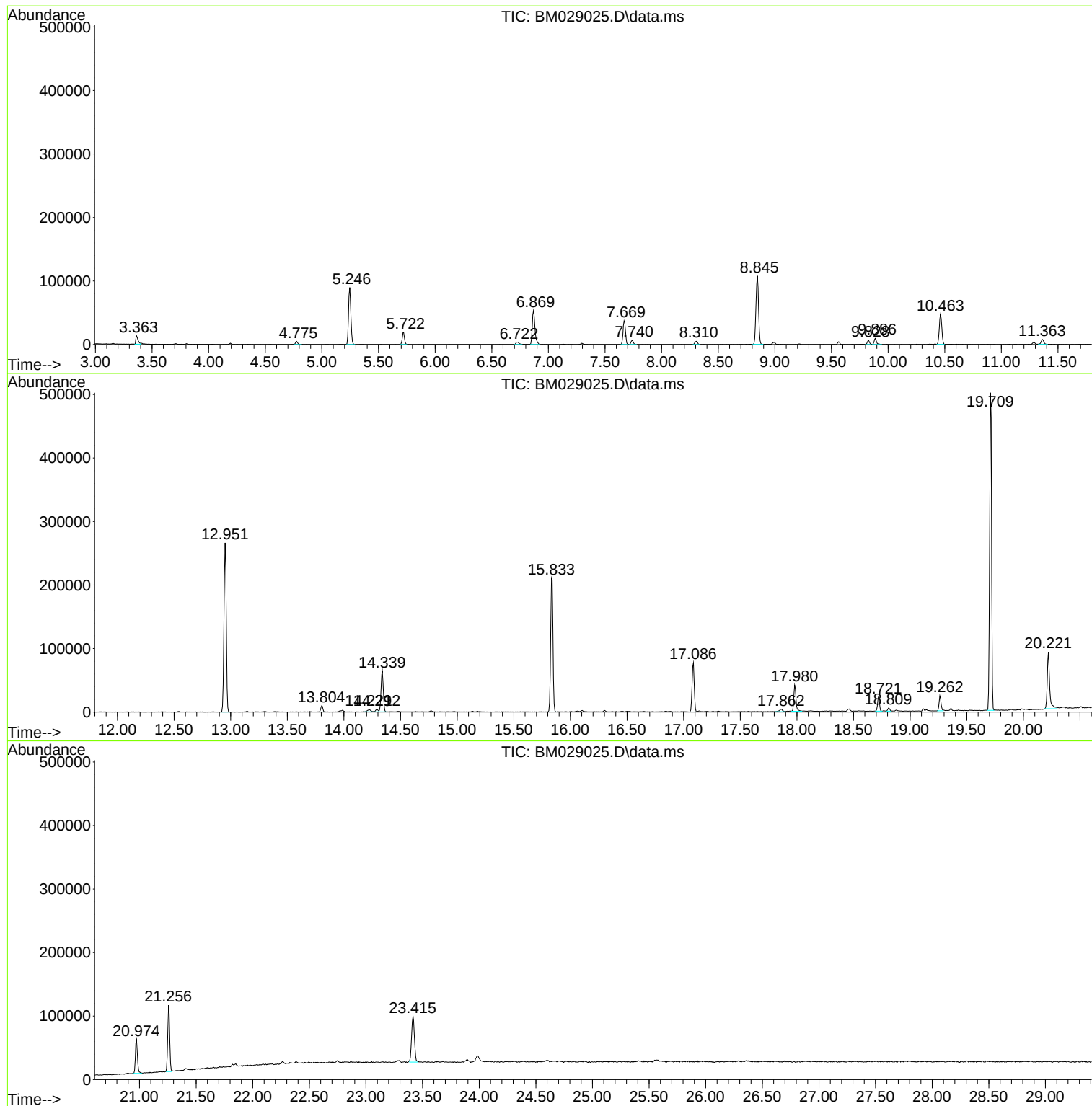
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ClientSampleId :
IDW-GW806-AQ-030321

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022321.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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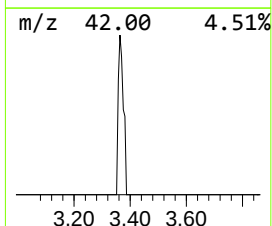
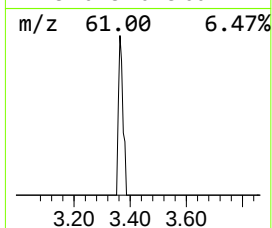
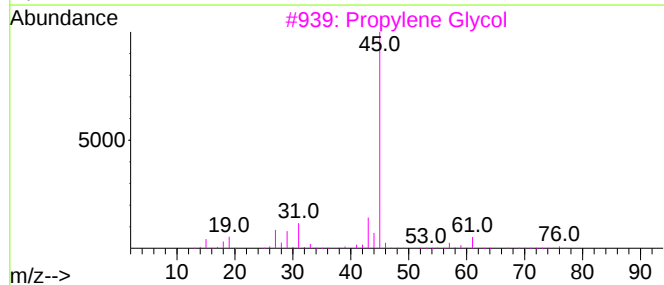
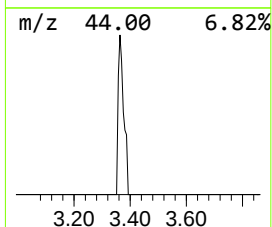
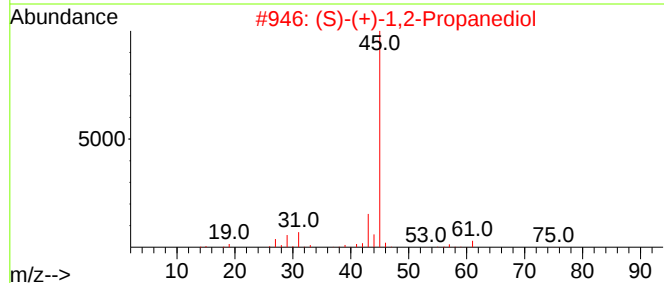
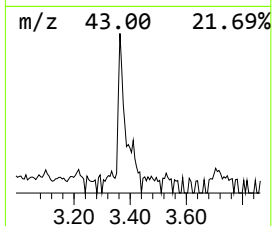
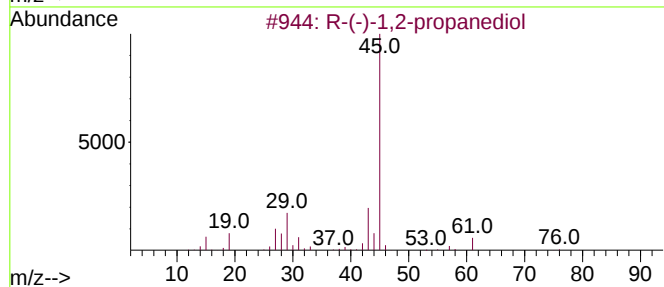
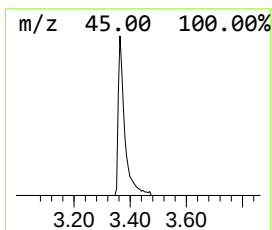
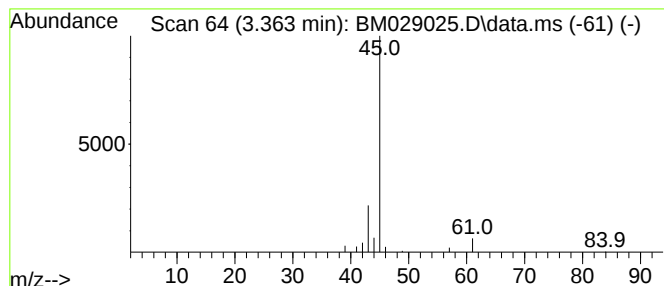
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 R-(-)-1,2-propanediol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.363	6.39 ng	17667	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	64
2			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	64
3			Propylene Glycol	76	C3H8O2	000057-55-6	9
4			Isopropyl Alcohol	60	C3H8O	000067-63-0	9
5			Formamide	45	CH3NO	000075-12-7	5



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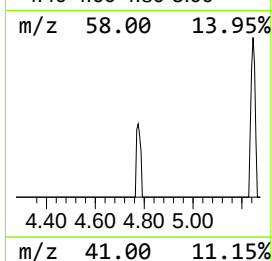
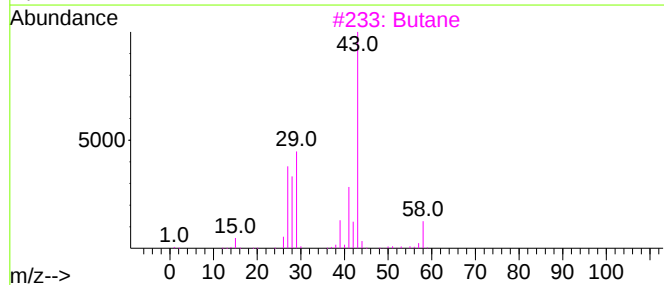
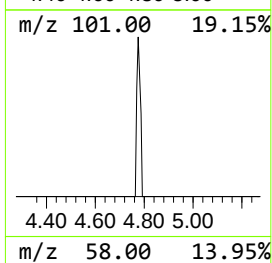
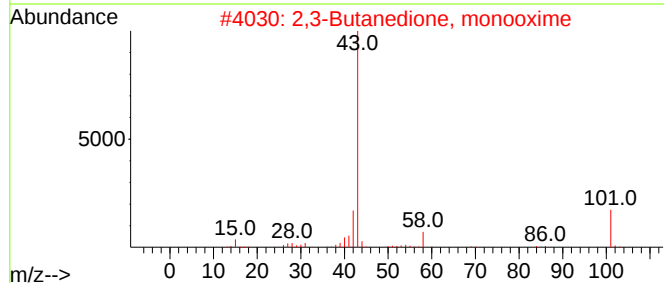
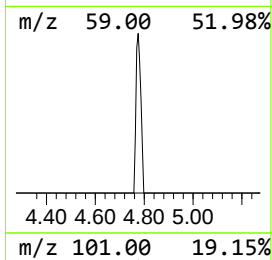
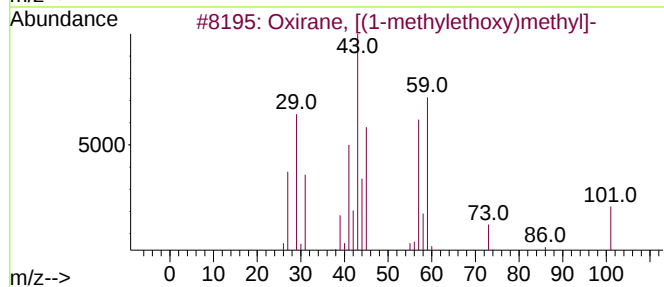
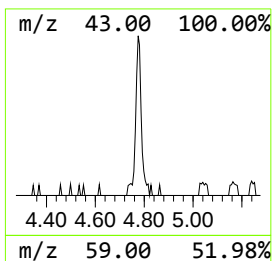
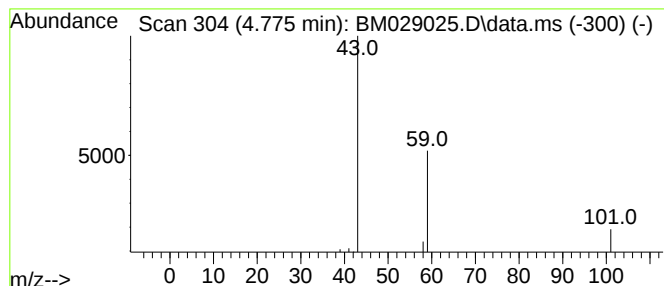
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022321.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown4.775 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.775	2.47 ng	6825	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	9
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	7
3			Butane	58	C4H10	000106-97-8	7
4			Diacetamide	101	C4H7NO2	000625-77-4	4
5			Propylamine	59	C3H9N	000107-10-8	4



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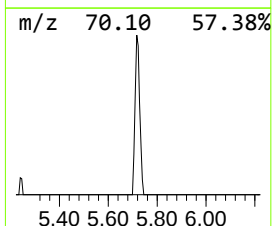
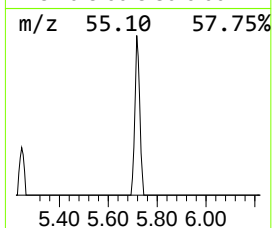
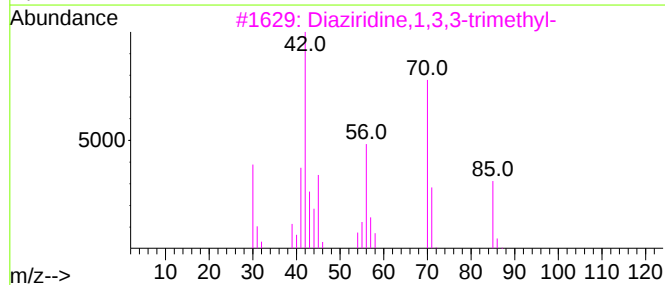
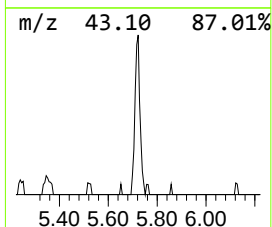
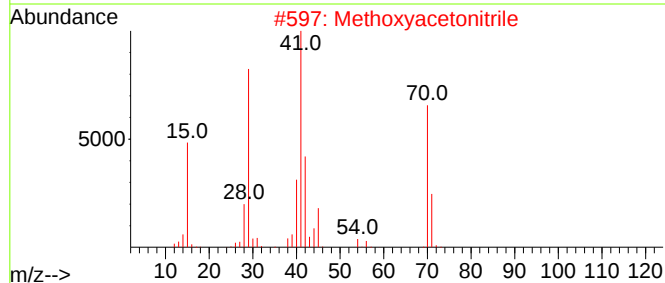
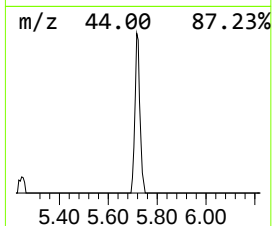
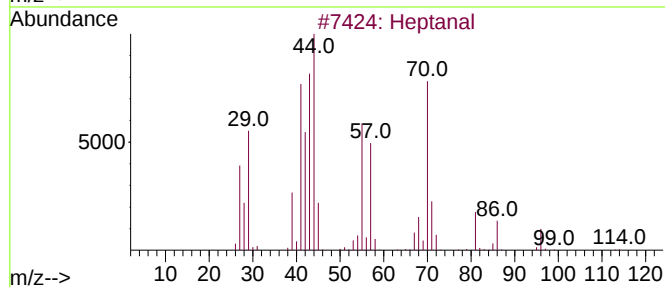
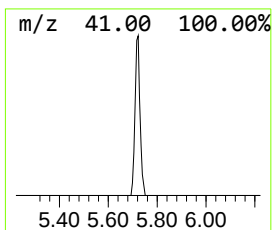
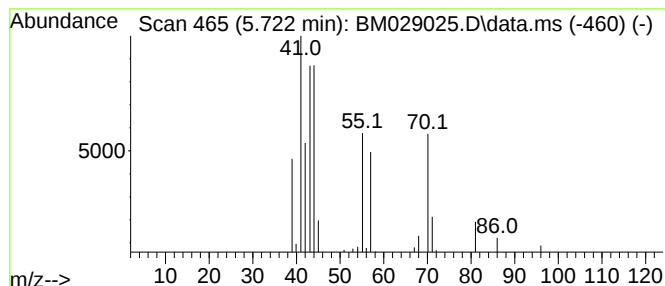
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Heptanal Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.722	9.38 ng	25926	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptanal	114	C7H14O	000111-71-7	83
2			Methoxyacetoneitrile	71	C3H5NO	001738-36-9	35
3			Diaziridine,1,3,3-trimethyl-	86	C4H10N2	040711-15-7	32
4			2-Pentene, (E)-	70	C5H10	000646-04-8	25
5			Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	25



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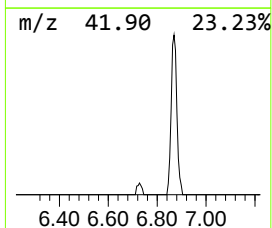
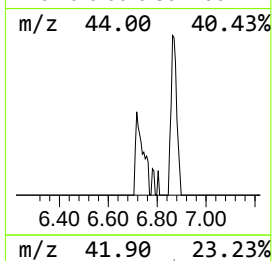
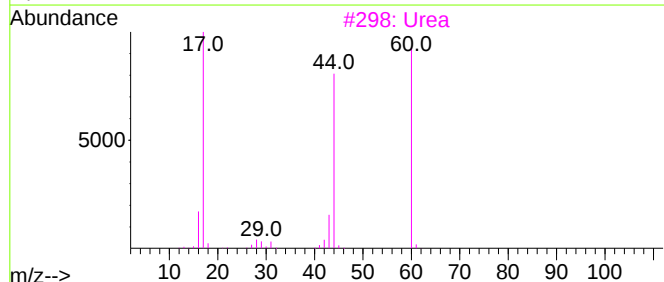
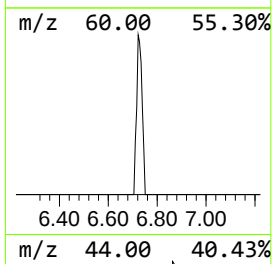
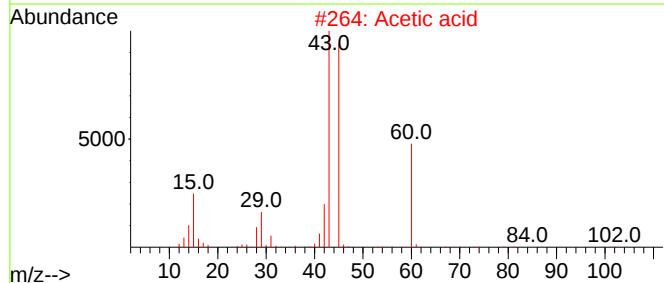
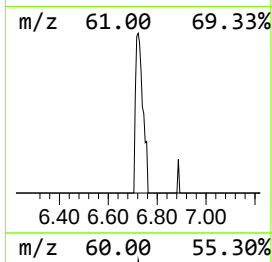
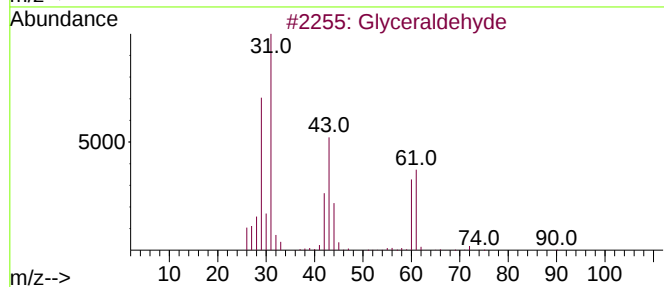
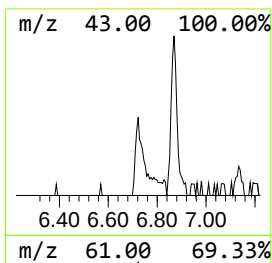
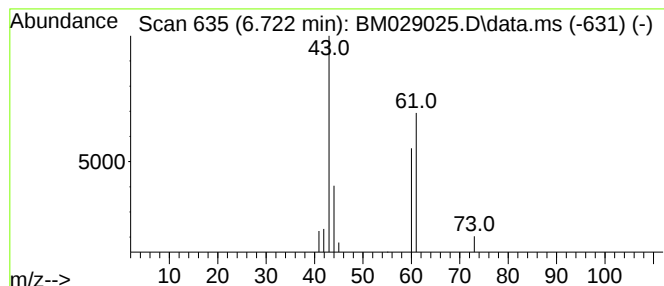
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown6.722 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.722	2.82 ng	7798	1,4-Dichlorobenzene-d4	7.669

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Glyceraldehyde	90	C3H6O3	000056-82-6	9
2		Acetic acid	60	C2H4O2	000064-19-7	7
3		Urea	60	CH4N2O	000057-13-6	5
4		Hydrazine, 1,1-dimethyl-	60	C2H8N2	000057-14-7	4
5		o-Ethylhydroxylamine	61	C2H7NO	1000322-42-5	4



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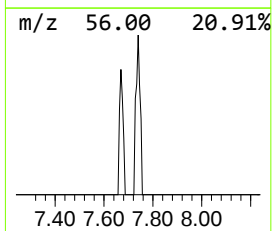
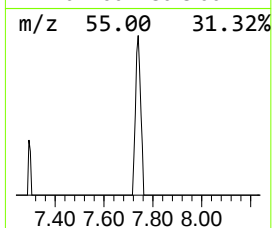
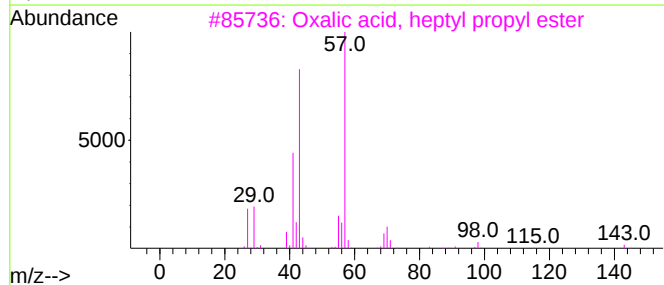
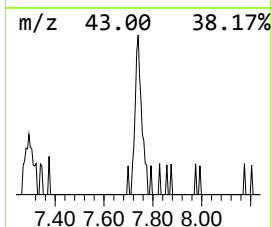
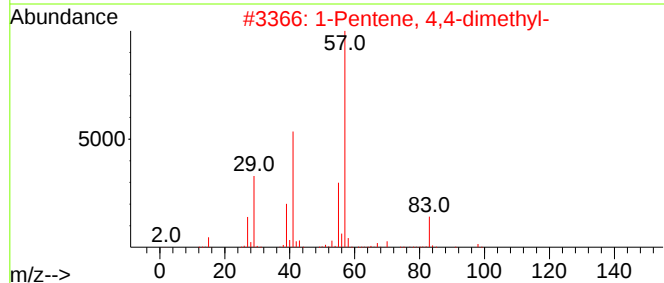
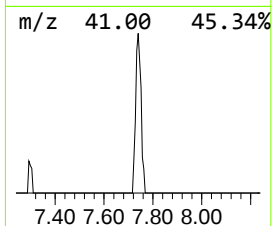
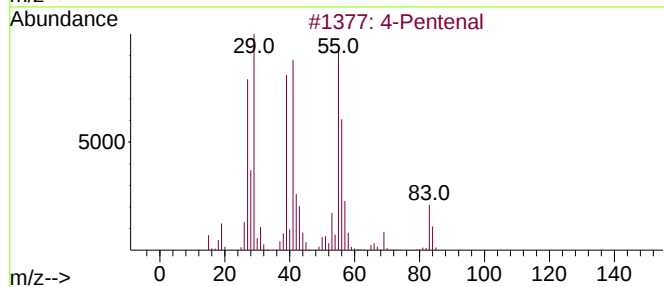
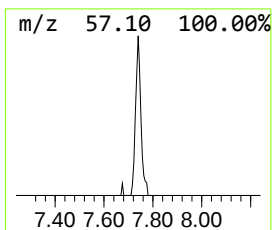
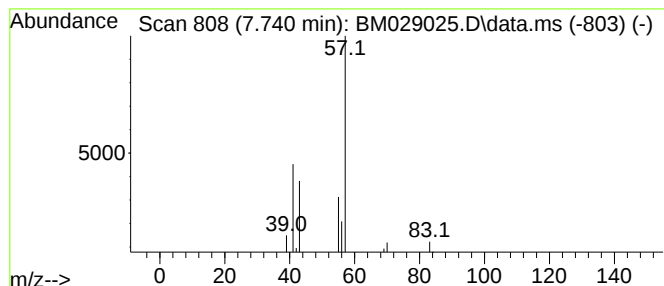
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 unknown7.740 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.740	3.54 ng	9795	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Pentenal	84	C5H8O	002100-17-6	23
2			1-Pentene, 4,4-dimethyl-	98	C7H14	000762-62-9	9
3			Oxalic acid, heptyl propyl ester	230	C12H22O4	1000309-26-0	9
4			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	4
5			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	4



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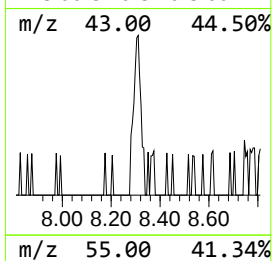
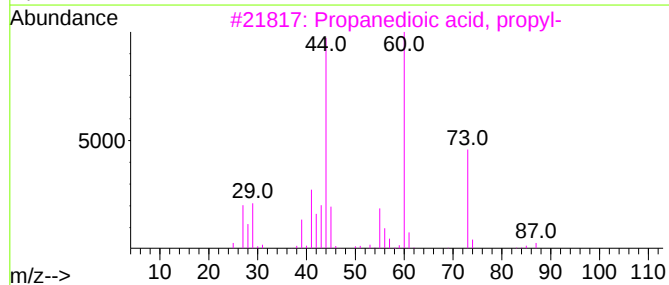
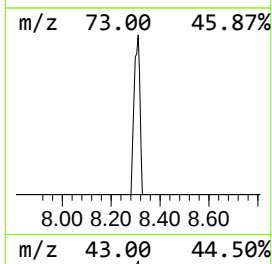
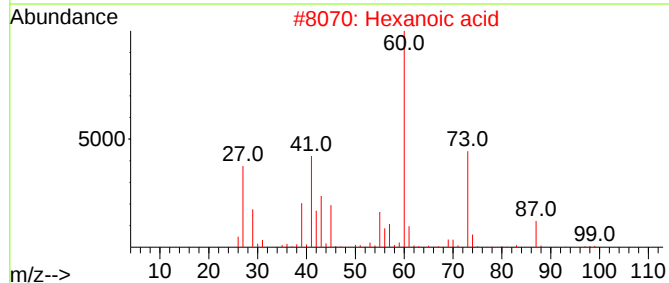
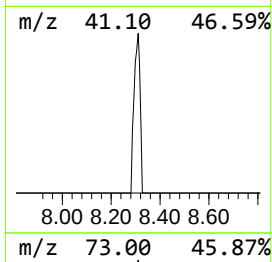
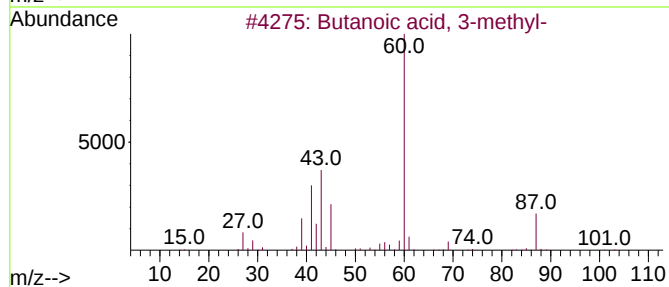
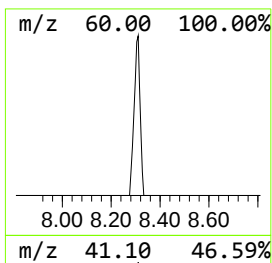
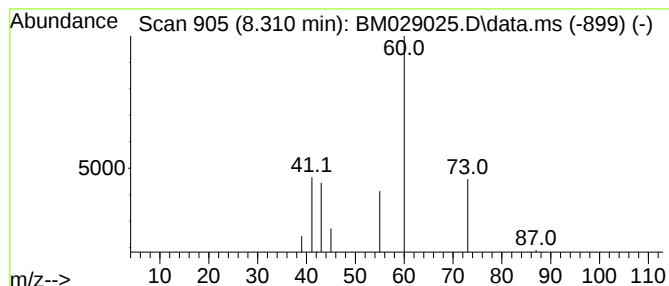
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown8.310 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.310	3.15 ng	8723	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	40
2			Hexanoic acid	116	C6H12O2	000142-62-1	40
3			Propanedioic acid, propyl-	146	C6H10O4	000616-62-6	9
4			Formic acid hydrazide	60	CH4N2O	000624-84-0	7
5			1-Pentanamine	87	C5H13N	000110-58-7	7



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
Data File : BM029025.D
Acq On : 08 Mar 2021 18:55
Operator : CG/JU
Sample : M1575-04
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
IDW-GW806-AQ-030321

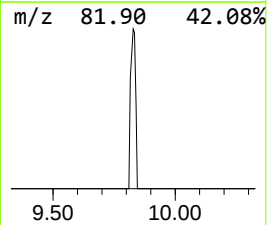
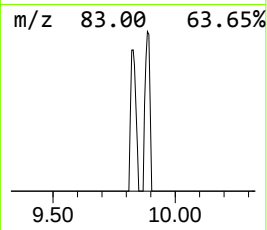
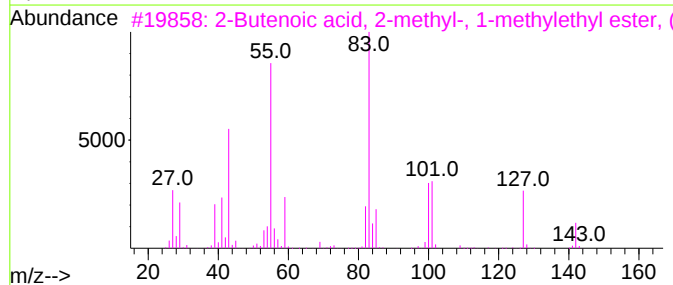
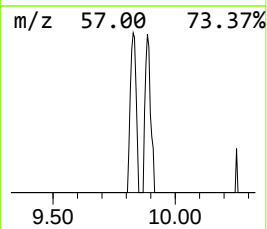
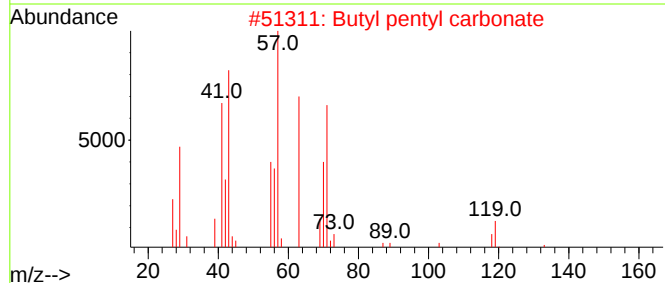
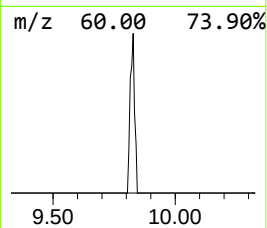
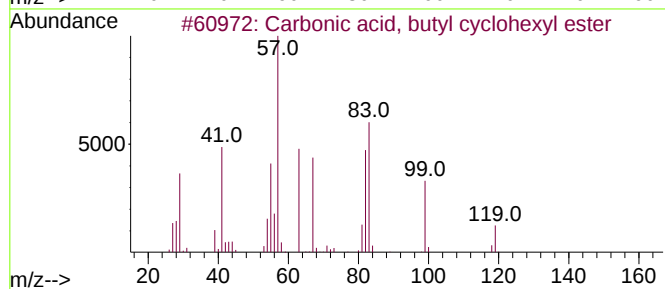
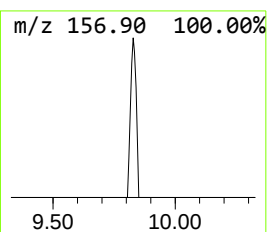
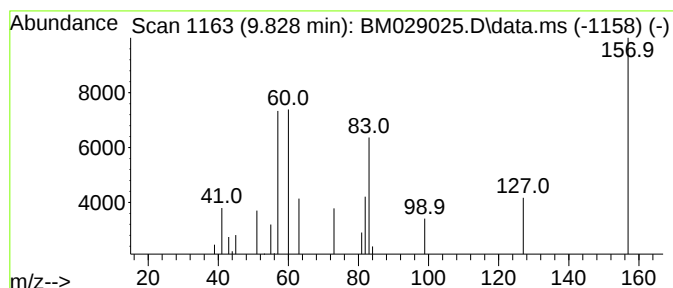
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown9.828 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.828	2.41 ng	8834	Naphthalene-d8	10.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, butyl cyclohexyl ...	200	C11H20O3	1000357-84-2	37
2			Butyl pentyl carbonate	188	C10H20O3	083073-03-4	9
3			2-Butenoic acid, 2-methyl-, 1-me...	142	C8H14O2	061692-76-0	9
4			Carbonic acid, isobutyl cyclohex...	200	C11H20O3	1000357-82-7	9
5			Octanoic acid, 2-ethylhexyl ester	256	C16H32O2	063321-70-0	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
Data File : BM029025.D
Acq On : 08 Mar 2021 18:55
Operator : CG/JU
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Misc :
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ClientSampleId :
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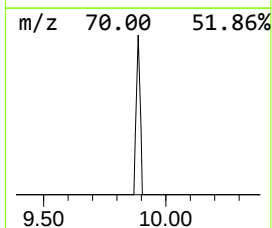
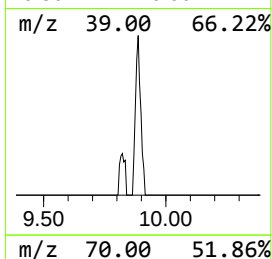
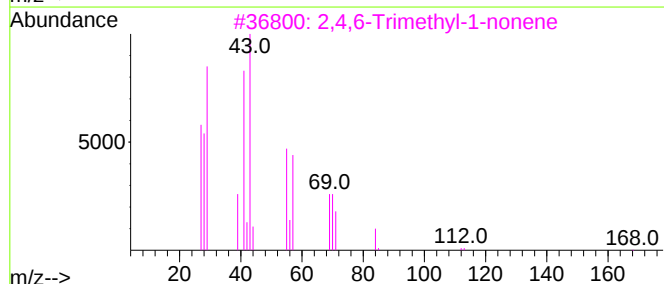
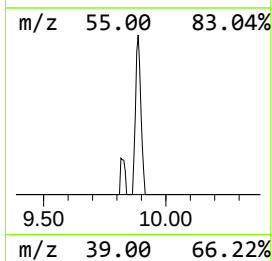
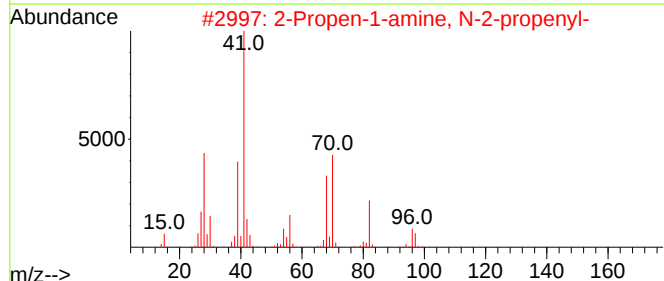
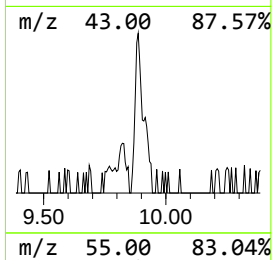
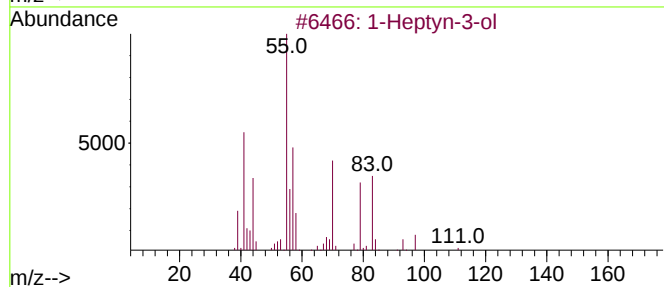
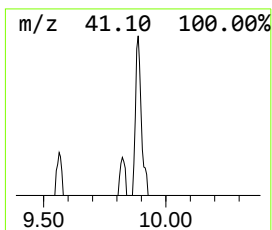
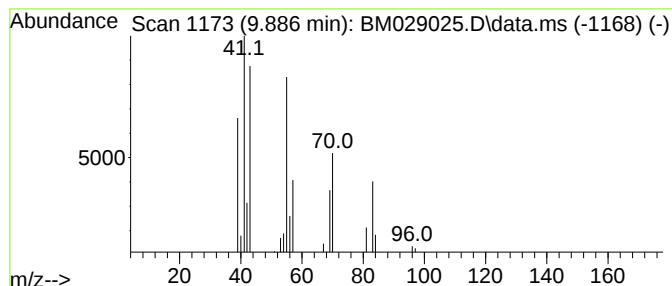
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1-Heptyn-3-ol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.886	3.94 ng	14456	Naphthalene-d8	10.463

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heptyn-3-ol	112	C7H12O	007383-19-9	50
2		2-Propen-1-amine, N-2-propenyl-	97	C6H11N	000124-02-7	47
3		2,4,6-Trimethyl-1-nonene	168	C12H24	055771-40-9	40
4		Allyl methallyl ether	112	C7H12O	014289-96-4	38
5		4-Pentenal	84	C5H8O	002100-17-6	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
Data File : BM029025.D
Acq On : 08 Mar 2021 18:55
Operator : CG/JU
Sample : M1575-04
Misc :
ALS Vial : 9 Sample Multiplier: 1

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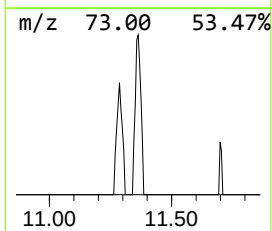
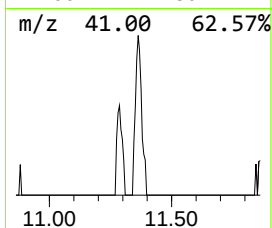
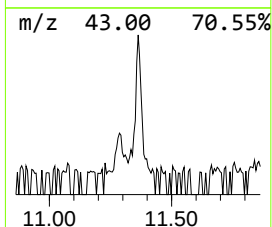
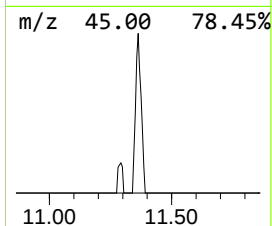
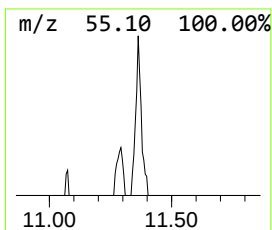
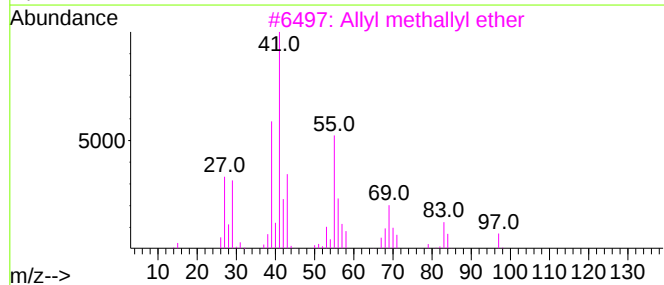
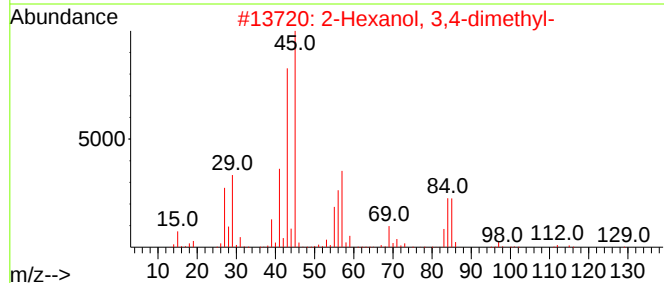
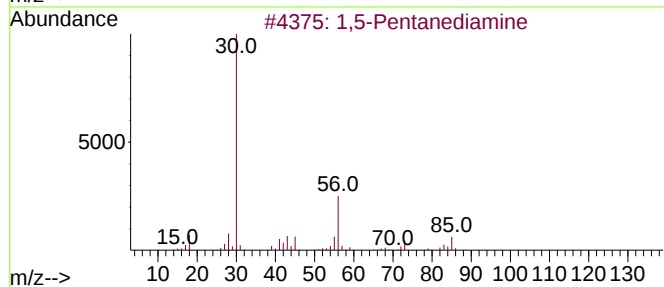
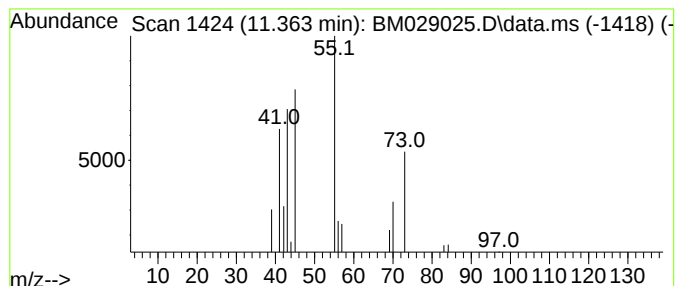
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown11.363 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.363	3.62 ng	13270	Naphthalene-d8	10.463

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,5-Pentanediamine	102	C5H14N2	000462-94-2	33
2		2-Hexanol, 3,4-dimethyl-	130	C8H18O	019550-05-1	25
3		Allyl methallyl ether	112	C7H12O	014289-96-4	22
4		3-Hexene, (Z)-	84	C6H12	007642-09-3	14
5		2-Octanol, (S)-	130	C8H18O	006169-06-8	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
Data File : BM029025.D
Acq On : 08 Mar 2021 18:55
Operator : CG/JU
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Misc :
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Instrument :
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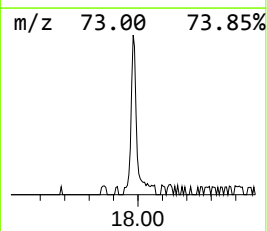
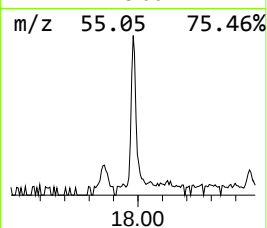
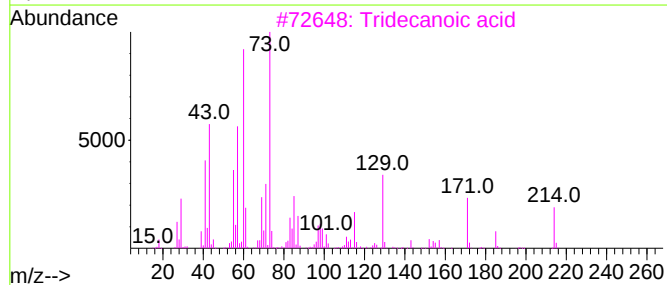
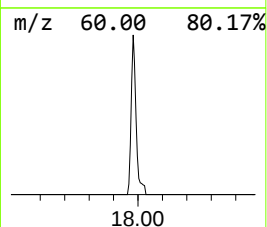
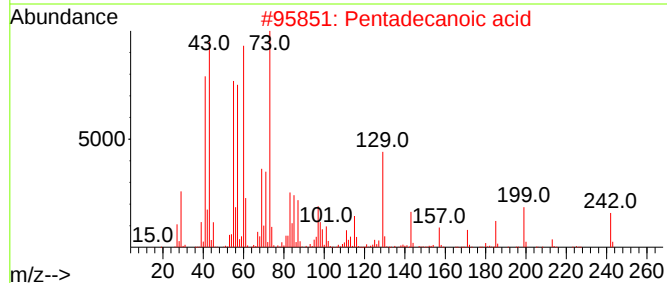
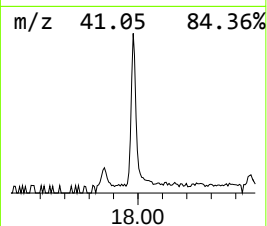
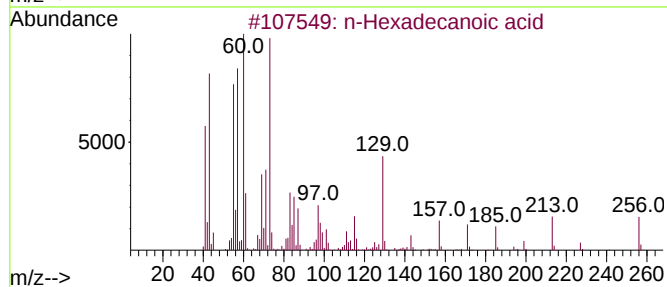
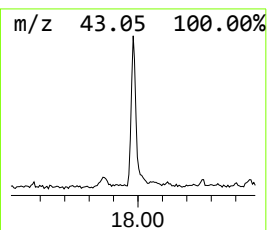
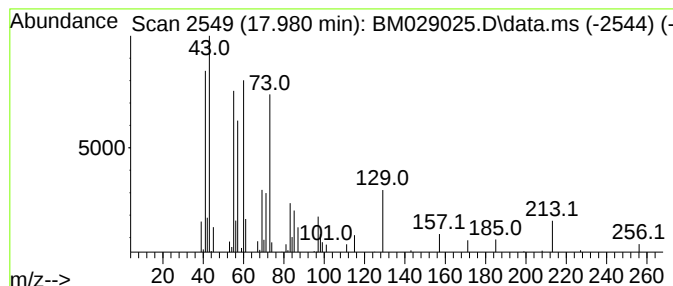
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.980	10.58 ng	54491	Phenanthrene-d10	17.086

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3			Tridecanoic acid	214	C13H26O2	000638-53-9	72
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	58
5			n-Decanoic acid	172	C10H20O2	000334-48-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
Data File : BM029025.D
Acq On : 08 Mar 2021 18:55
Operator : CG/JU
Sample : M1575-04
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
IDW-GW806-AQ-030321

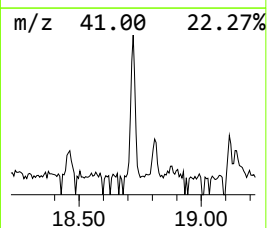
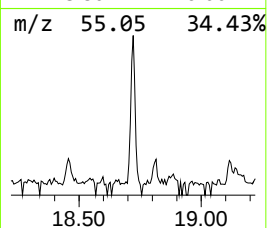
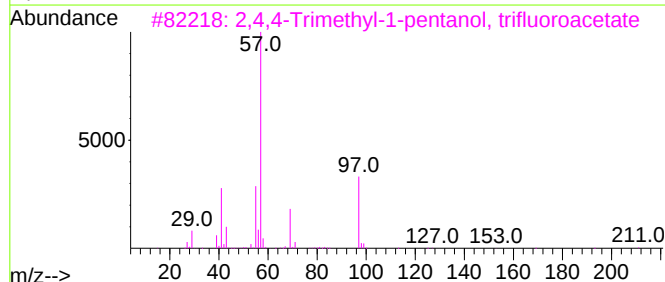
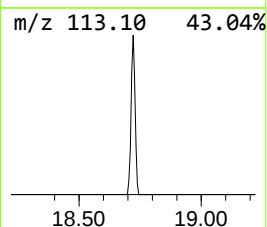
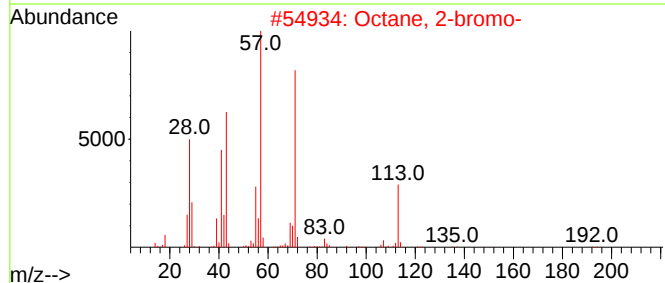
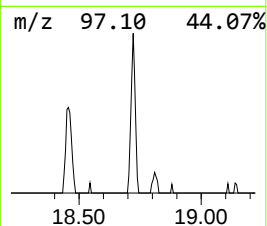
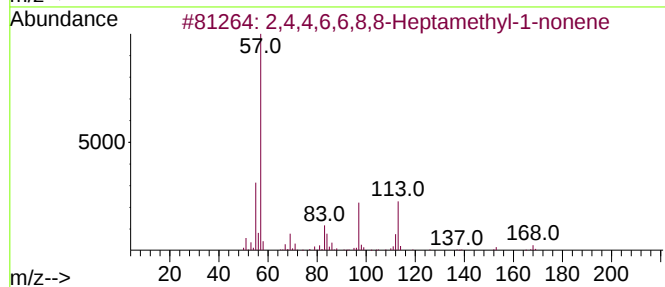
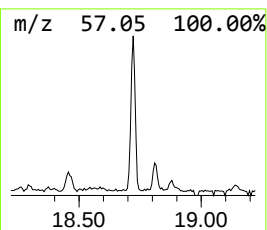
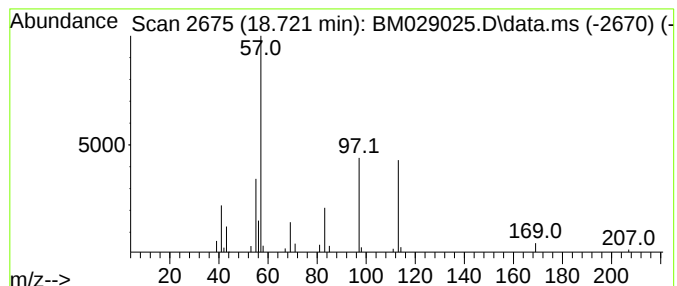
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2,4,4,6,6,8,8-Heptamethyl-1... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.721	5.21 ng	26825	Phenanthrene-d10	17.086

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	56	
2	Octane, 2-bromo-	192	C8H17Br	000557-35-7	37	
3	2,4,4-Trimethyl-1-pentanol, trif...	226	C10H17F3O2	1000365-19-5	35	
4	Pentane, 2,2,4-trimethyl-4-nitro-	159	C8H17NO2	005342-78-9	25	
5	Cyclohexane, 1,1'-[1,2-bis(1,1-d...	306	C22H42	065149-85-1	25	



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
Data File : BM029025.D
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Misc :
ALS Vial : 9 Sample Multiplier: 1

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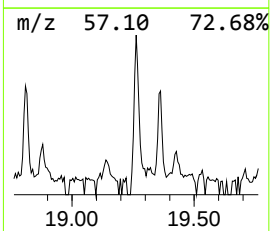
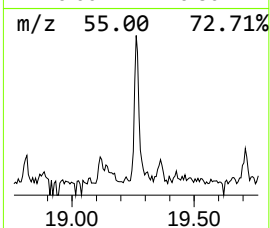
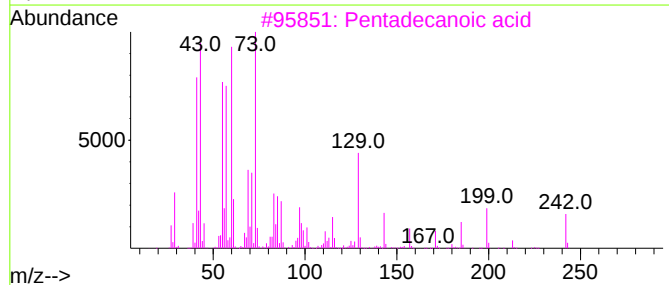
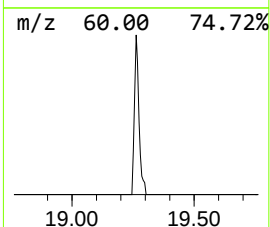
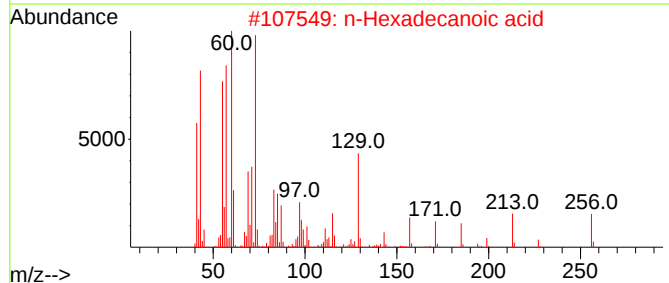
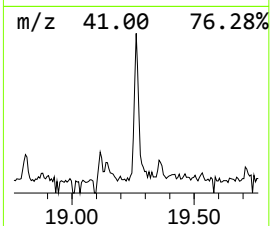
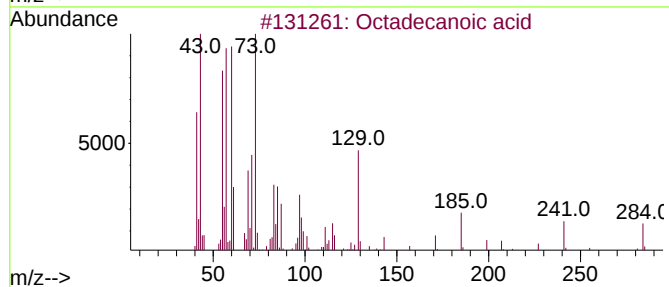
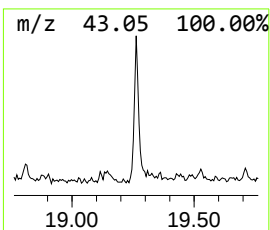
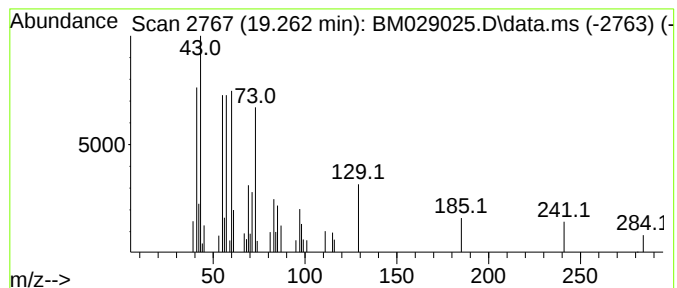
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.262	4.59 ng	28873	Chrysene-d12	21.256

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	96
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	74
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	64
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	59
5			Tridecanoic acid	214	C13H26O2	000638-53-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
Data File : BM029025.D
Acq On : 08 Mar 2021 18:55
Operator : CG/JU
Sample : M1575-04
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
IDW-GW806-AQ-030321

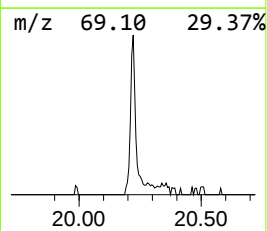
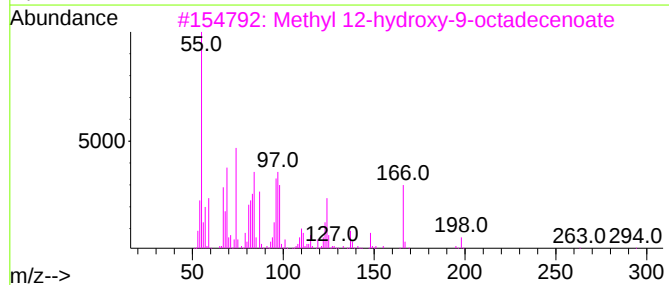
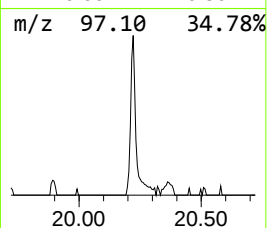
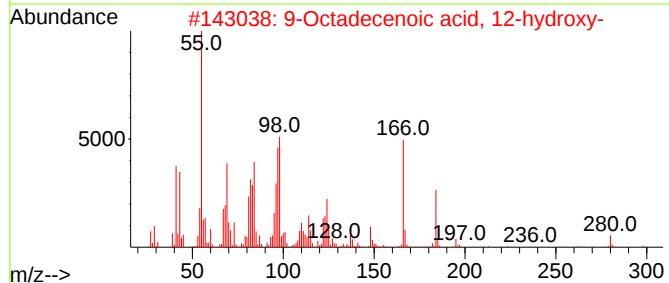
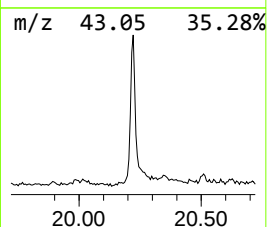
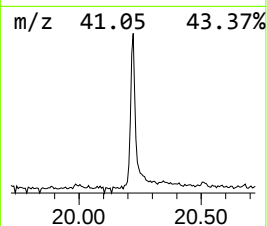
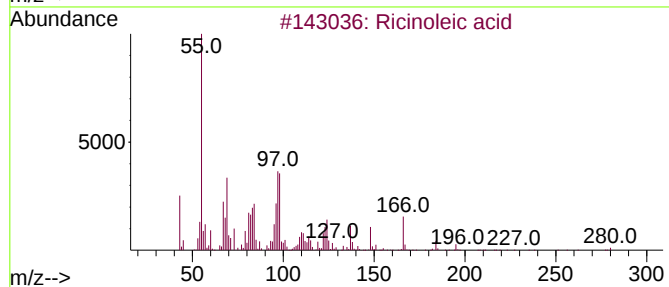
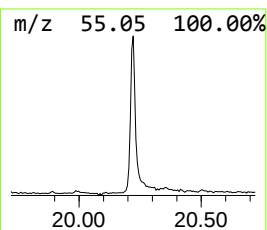
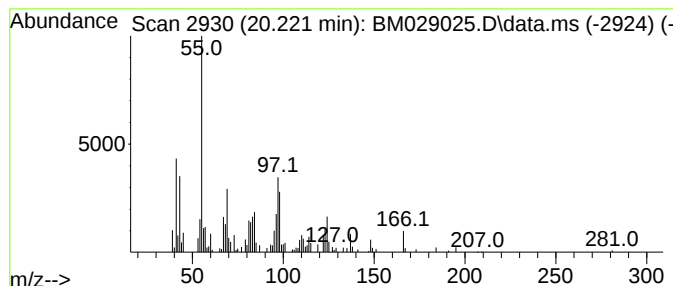
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Ricinoleic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.221	20.72 ng	130246	Chrysene-d12	21.256

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ricinoleic acid	298	C18H34O3	000141-22-0	91
2			9-Octadecenoic acid, 12-hydroxy-	298	C18H34O3	007431-95-0	72
3			Methyl 12-hydroxy-9-octadecenoate	312	C19H36O3	1000336-28-8	53
4			E-9-Tetradecenoic acid	226	C14H26O2	1000131-35-8	38
5			1-Nonylcycloheptane	224	C16H32	1000371-47-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
Data File : BM029025.D
Acq On : 08 Mar 2021 18:55
Operator : CG/JU
Sample : M1575-04
Misc :
ALS Vial : 9 Sample Multiplier: 1

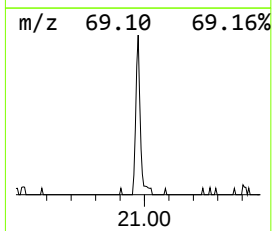
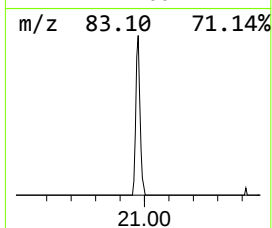
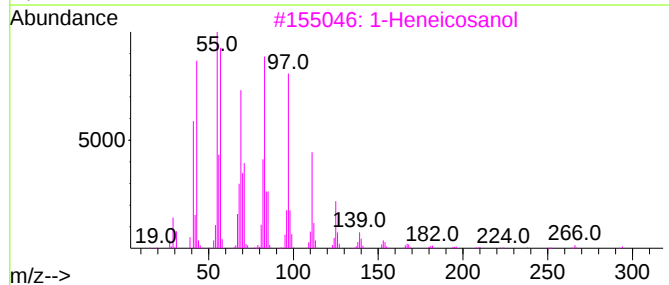
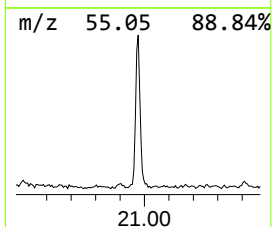
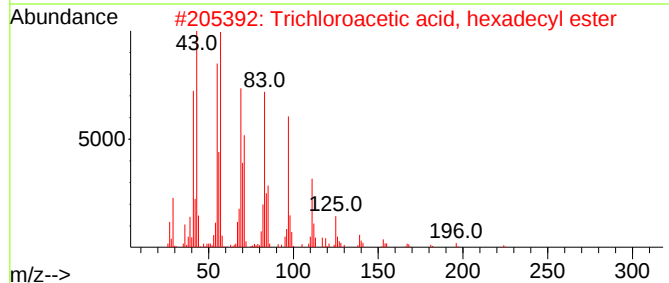
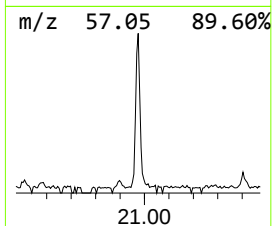
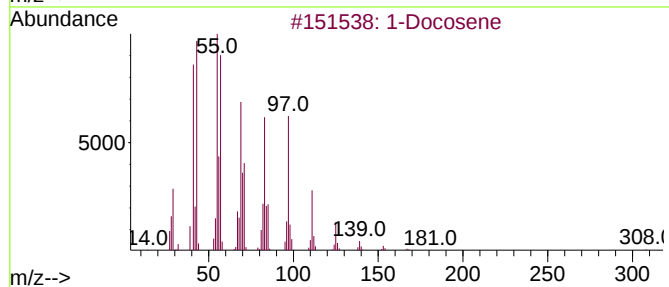
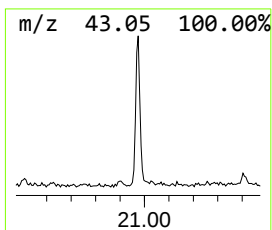
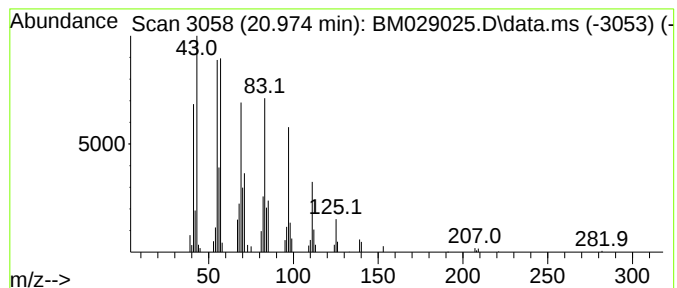
Instrument :
BNA_M
ClientSampleId :
IDW-GW806-AQ-030321

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1-Docosene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.974	10.77 ng	67676	Chrysene-d12		21.256	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Docosene		308	C22H44	001599-67-3	95
2	Trichloroacetic acid, hexadecyl ...		386	C18H33Cl3O2	074339-54-1	95
3	1-Heneicosanol		312	C21H44O	015594-90-8	95
4	Trichloroacetic acid, pentadecyl...		372	C17H31Cl3O2	074339-53-0	94
5	Heptadecyl trifluoroacetate		352	C19H35F3O2	1000351-87-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
 Data File : BM029025.D
 Acq On : 08 Mar 2021 18:55
 Operator : CG/JU
 Sample : M1575-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 IDW-GW806-AQ-030321

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022321.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
R-(-)-1,2-propa...	3.363	6.4	ng	17667	1	7.669	55302	20.0
unknown4.775	4.775	2.5	ng	6825	1	7.669	55302	20.0
Heptanal	5.722	9.4	ng	25926	1	7.669	55302	20.0
unknown6.722	6.722	2.8	ng	7798	1	7.669	55302	20.0
unknown7.740	7.740	3.5	ng	9795	1	7.669	55302	20.0
unknown8.310	8.310	3.1	ng	8723	1	7.669	55302	20.0
unknown9.828	9.828	2.4	ng	8834	2	10.463	73384	20.0
1-Heptyn-3-ol	9.886	3.9	ng	14456	2	10.463	73384	20.0
unknown11.363	11.363	3.6	ng	13270	2	10.463	73384	20.0
n-Hexadecanoic ...	17.980	10.6	ng	54491	4	17.086	103022	20.0
2,4,4,6,6,8,8-H...	18.721	5.2	ng	26825	4	17.086	103022	20.0
Octadecanoic acid	19.262	4.6	ng	28873	5	21.256	125715	20.0
Ricinoleic acid	20.221	20.7	ng	130246	5	21.256	125715	20.0
1-Docosene	20.974	10.8	ng	67676	5	21.256	125715	20.0