

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
 Data File : BM029031.D
 Acq On : 08 Mar 2021 22:34
 Operator : CG/JU
 Sample : M1582-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 VEOLIA-VNJ-208

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: BM029031.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.399	60	70	87	rBV	337171	722818	70.21%	10.726%
2	5.246	379	384	394	rBB	170410	236427	22.96%	3.508%
3	6.869	654	660	671	rBV	159752	246059	23.90%	3.651%
4	7.669	791	796	802	rVB	42335	59977	5.83%	0.890%
5	8.845	989	996	1004	rBV	94328	152341	14.80%	2.261%
6	10.463	1264	1271	1278	rBB	49520	77869	7.56%	1.156%
7	12.951	1688	1694	1702	rBB	206958	281630	27.35%	4.179%
8	13.310	1746	1755	1769	rBV	70954	160526	15.59%	2.382%
9	14.227	1900	1911	1915	rBV3	4773	11361	1.10%	0.169%
10	14.274	1915	1919	1925	rVV	17246	29328	2.85%	0.435%
11	14.339	1925	1930	1937	rVB	66878	95567	9.28%	1.418%
12	15.833	2178	2184	2191	rBV	166314	220300	21.40%	3.269%
13	16.757	2334	2341	2352	rBV	69288	102130	9.92%	1.516%
14	17.086	2391	2397	2401	rBV	76727	105836	10.28%	1.571%
15	17.980	2541	2549	2557	rBV2	30524	42095	4.09%	0.625%
16	18.751	2675	2680	2690	rBV	207523	280233	27.22%	4.159%
17	19.139	2742	2746	2753	rVB	22355	27380	2.66%	0.406%
18	19.262	2763	2767	2771	rBV	12799	15152	1.47%	0.225%
19	19.503	2800	2808	2816	rVB	20589	29521	2.87%	0.438%
20	19.709	2838	2843	2848	rVB	338421	374502	36.37%	5.557%
21	20.327	2939	2948	2959	rBV	520473	691713	67.18%	10.265%
22	20.974	3051	3058	3065	rBV2	40308	60703	5.90%	0.901%
23	21.174	3088	3092	3096	rBV	27003	28113	2.73%	0.417%
24	21.256	3100	3106	3110	rVV	114070	150638	14.63%	2.235%
25	21.292	3110	3112	3118	rVB	12285	15059	1.46%	0.223%
26	21.603	3159	3165	3187	rBV	763455	1029579	100.00%	15.279%
27	21.844	3203	3206	3211	rVB5	12062	16064	1.56%	0.238%
28	22.903	3379	3386	3410	rBV	493904	886112	86.07%	13.150%
29	23.415	3468	3473	3480	rVB2	75165	125926	12.23%	1.869%
30	23.985	3565	3570	3576	rBV9	7743	17046	1.66%	0.253%
31	24.644	3674	3682	3698	rBV	176049	446727	43.39%	6.629%

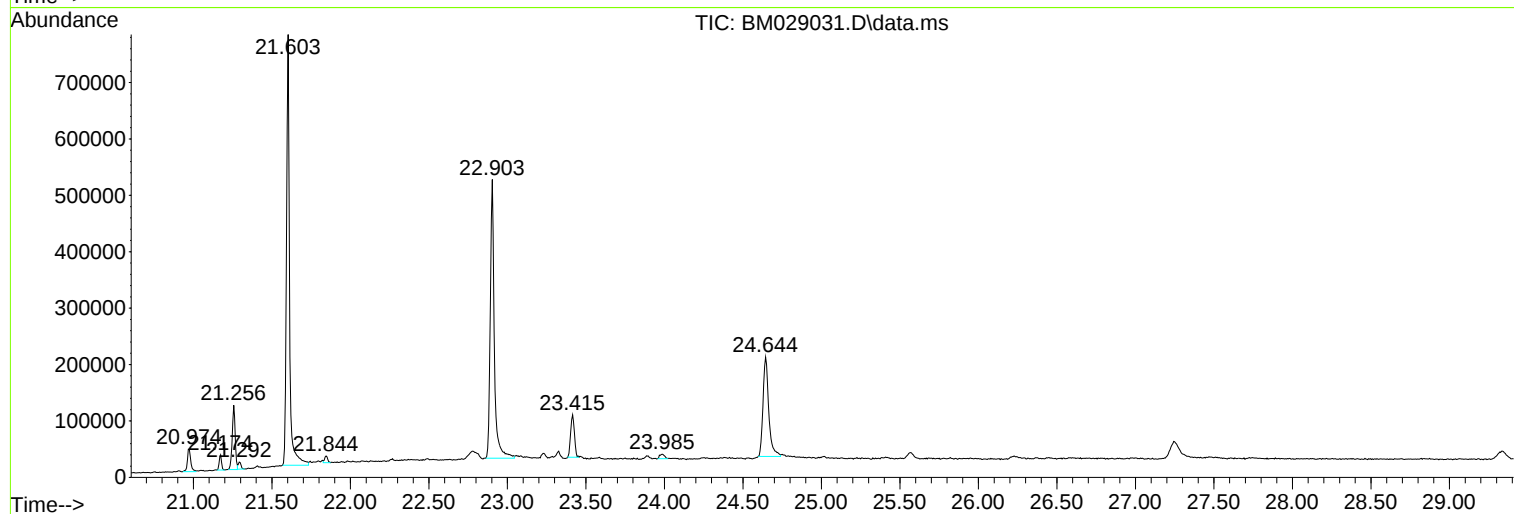
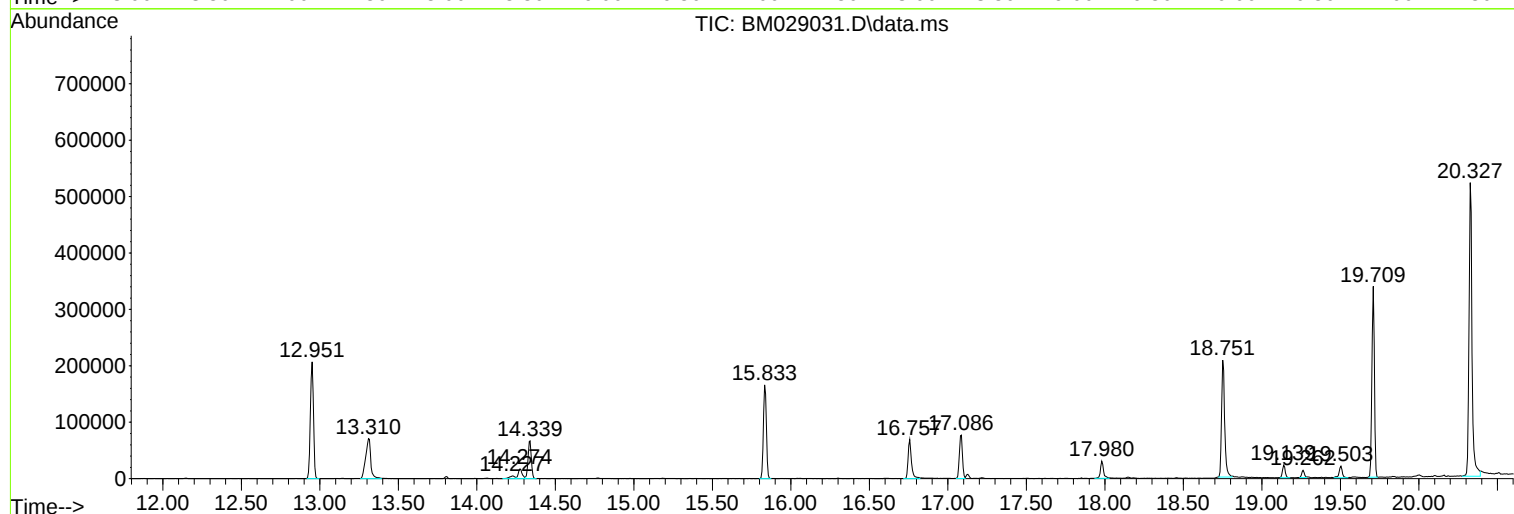
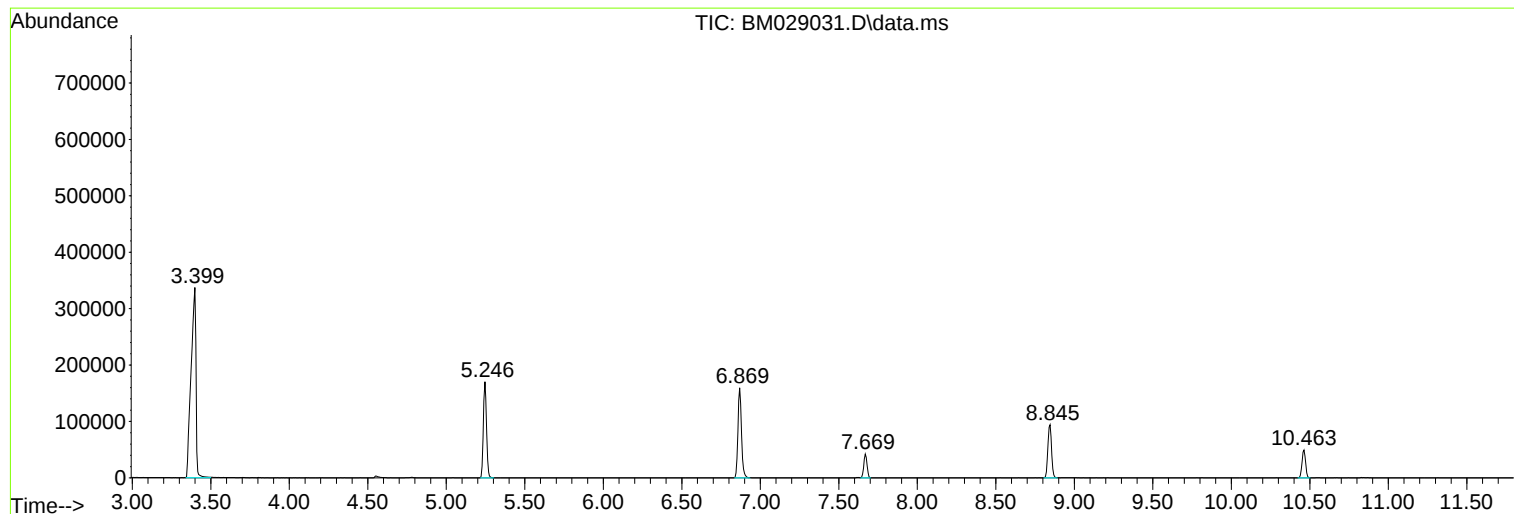
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Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
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ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
VEOLIA-VNJ-208

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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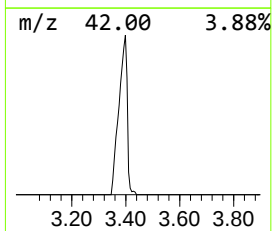
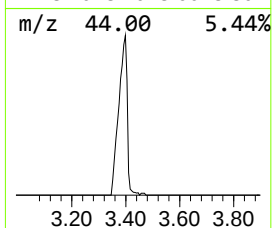
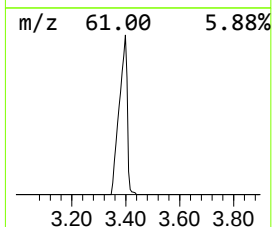
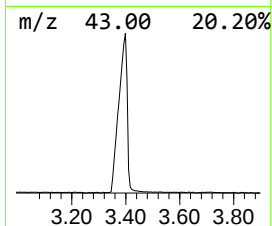
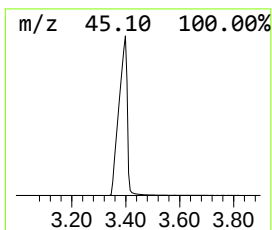
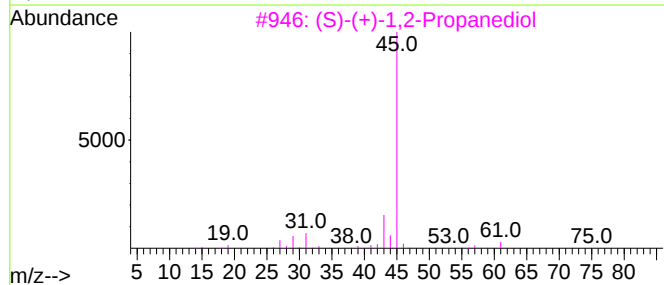
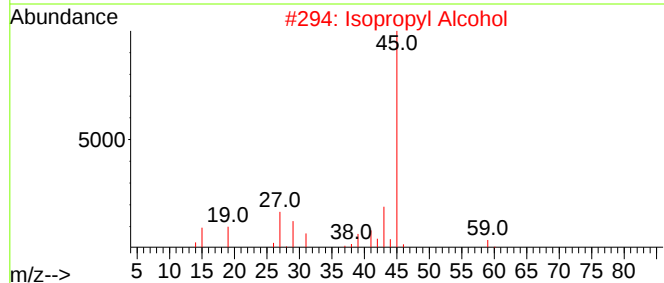
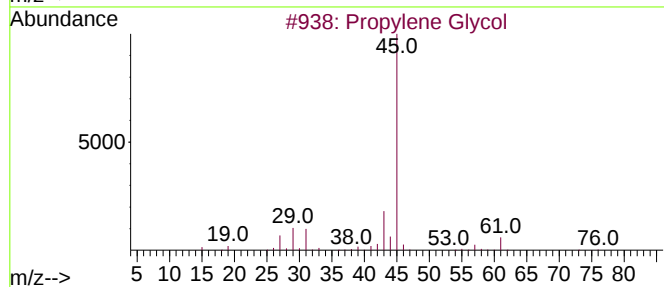
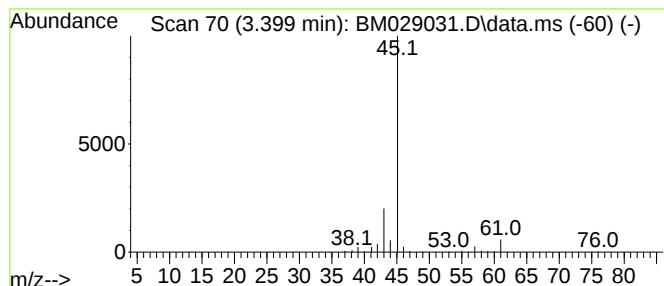
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.399	241.03 ng	722818	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	90
2			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
3			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	40
4			(R)-(-)-3-Methyl-2-butanol	88	C5H12O	001572-93-6	9
5			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	9



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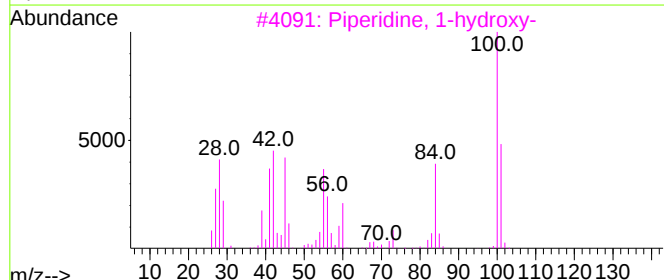
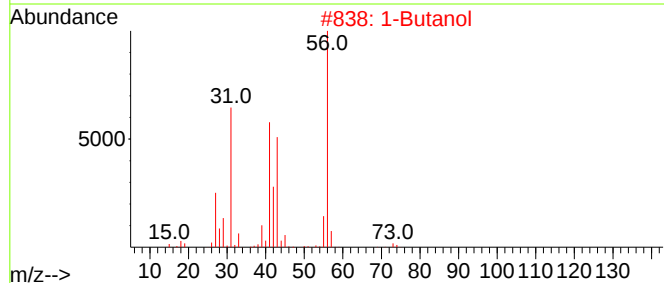
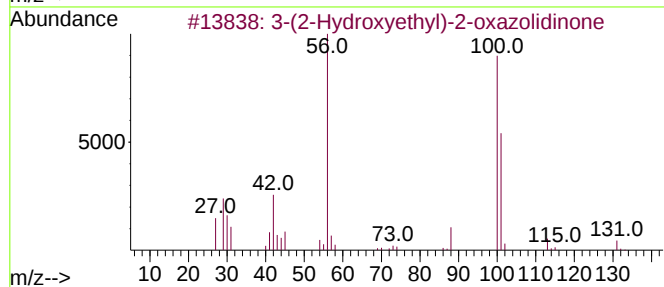
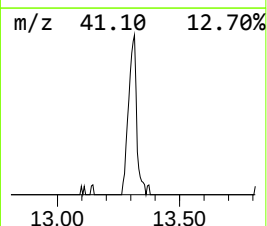
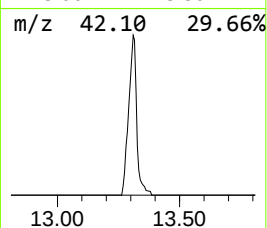
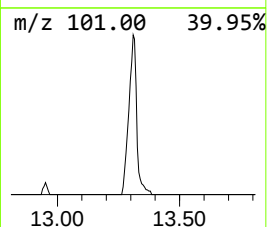
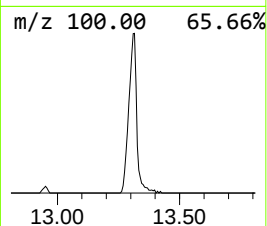
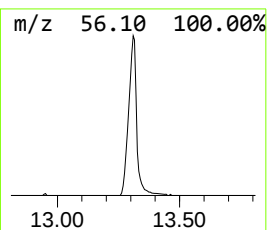
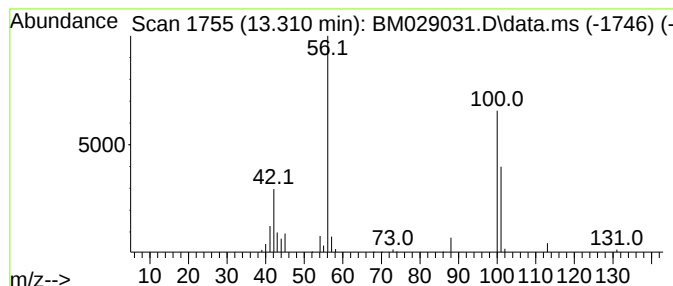
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 3-(2-Hydroxyethyl)-2-oxazol... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.310	33.59 ng	160526	Acenaphthene-d10	14.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-(2-Hydroxyethyl)-2-oxazolidinone	131	C5H9NO3	003356-88-5	78
2			1-Butanol	74	C4H10O	000071-36-3	43
3			Piperidine, 1-hydroxy-	101	C5H11NO	004801-58-5	37
4			Methanamine, N-(1-methylbutylidene...	99	C6H13N	022431-09-0	17
5			2H-1,4-Oxazine-4-propanoic acid,...	173	C8H15NO3	1000362-52-8	12



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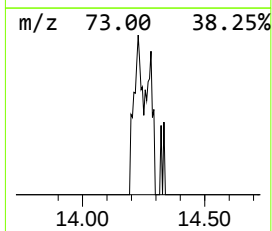
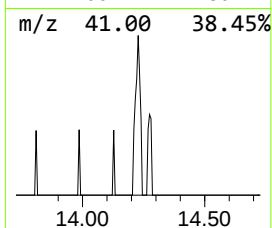
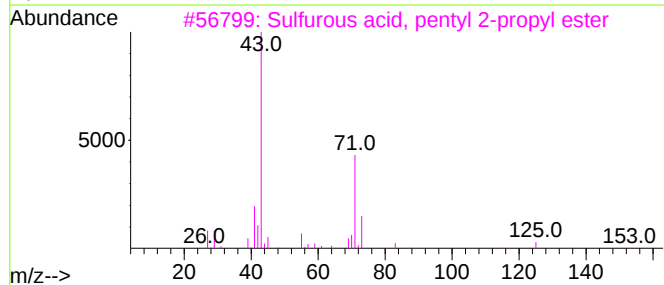
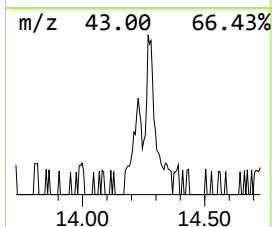
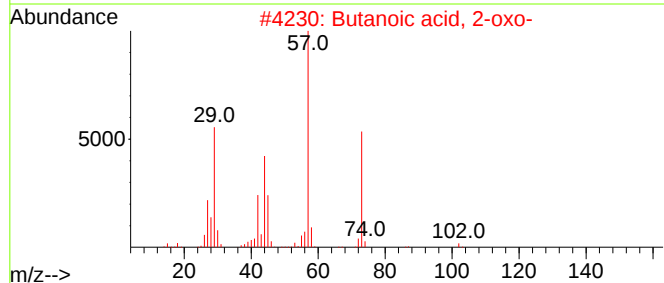
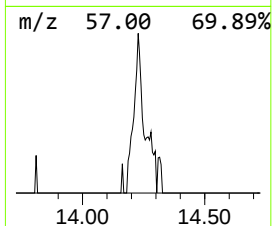
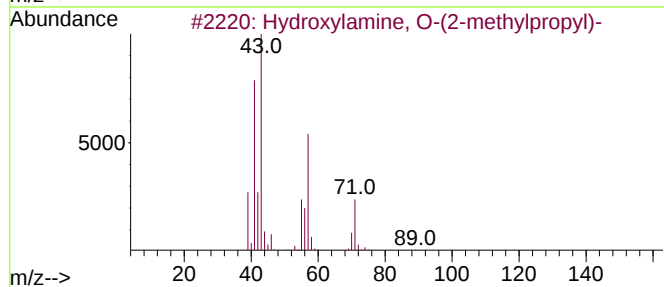
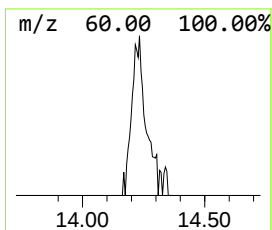
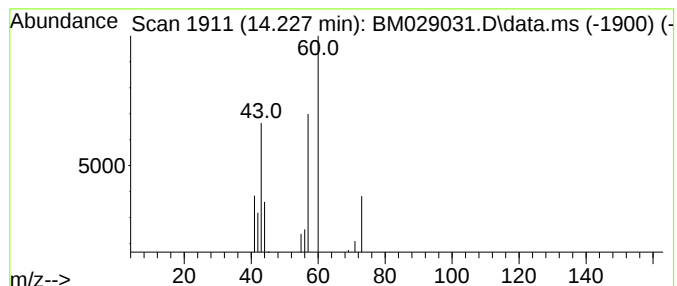
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown14.227 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.227	2.38 ng	11361	Acenaphthene-d10	14.339

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hydroxylamine, O-(2-methylpropyl)-	89	C4H11NO	005618-62-2	12
2		Butanoic acid, 2-oxo-	102	C4H6O3	000600-18-0	9
3		Sulfurous acid, pentyl 2-propyl ...	194	C8H18O3S	1000309-11-5	9
4		Azetidine, 2-methyl-	71	C4H9N	019812-49-8	7
5		2-Propenamide	71	C3H5NO	000079-06-1	7



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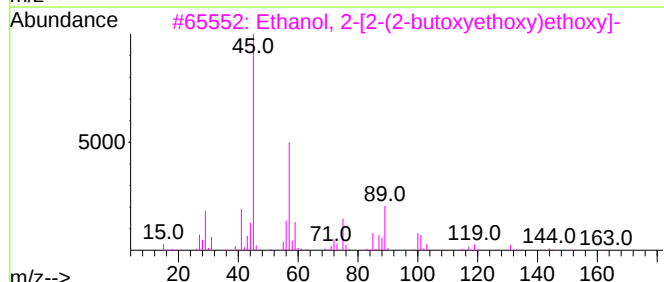
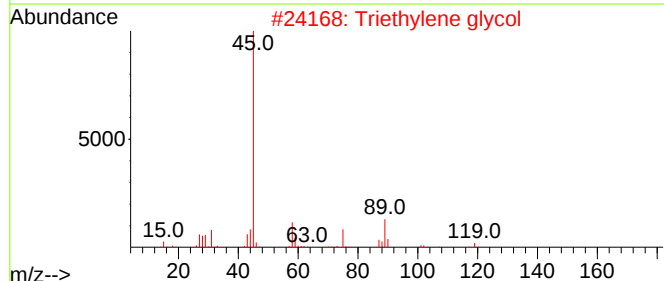
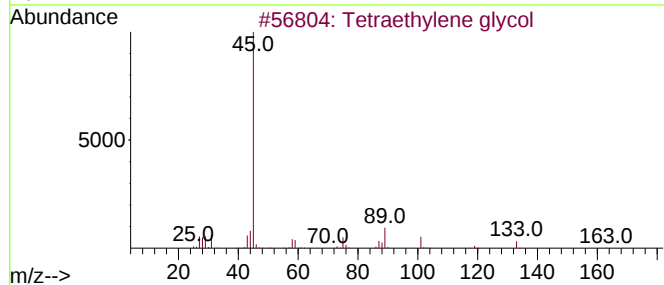
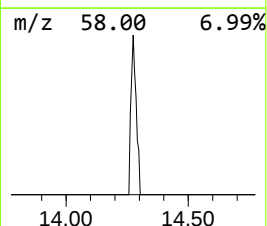
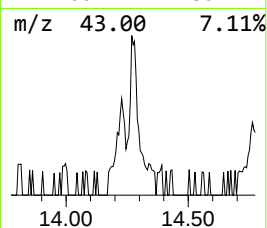
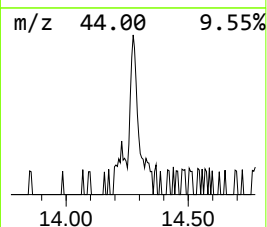
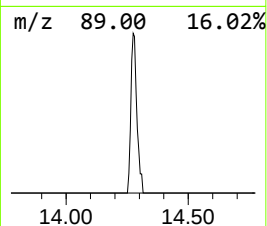
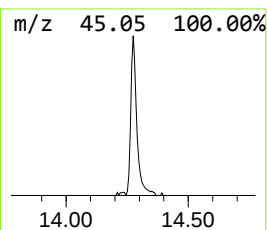
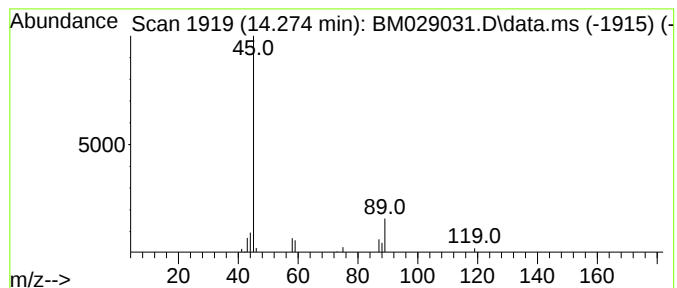
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Tetraethylene glycol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.274	6.14 ng	29328	Acenaphthene-d10	14.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetraethylene glycol	194	C8H18O5	000112-60-7	83
2			Triethylene glycol	150	C6H14O4	000112-27-6	78
3			Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	64
4			Ethanol, 2-[2-(2-methoxyethoxy)e...	164	C7H16O4	000112-35-6	64
5			Ethanol, 2-[2-(ethoxyloxy)ethoxy]-	132	C6H12O3	000929-37-3	56



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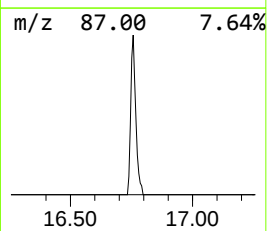
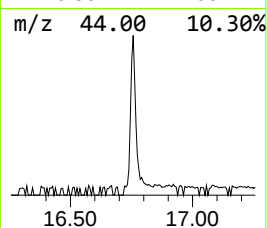
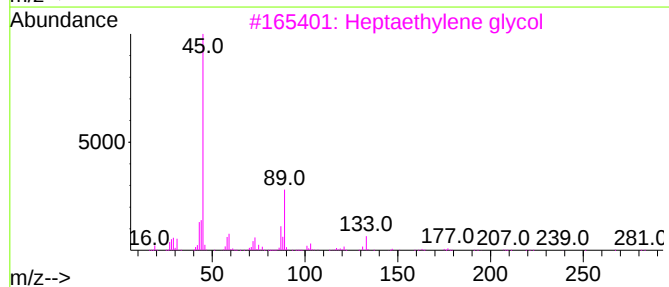
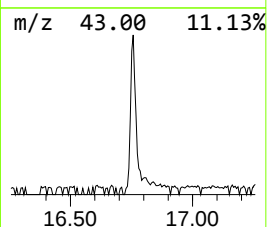
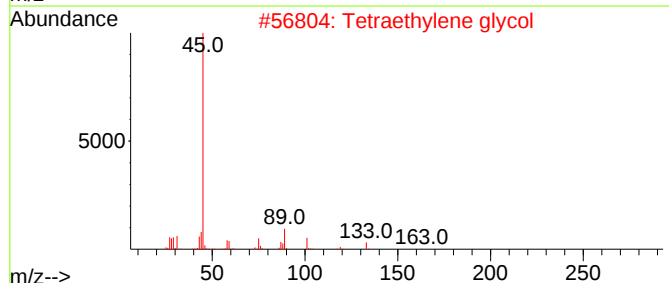
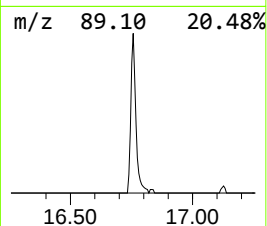
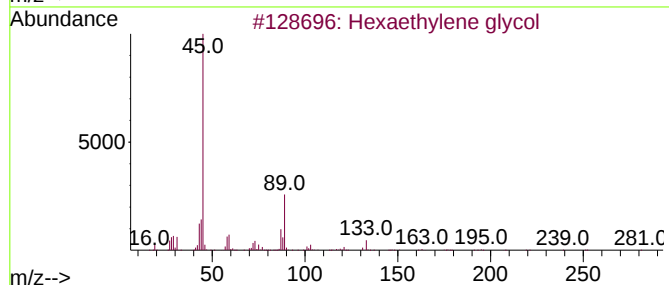
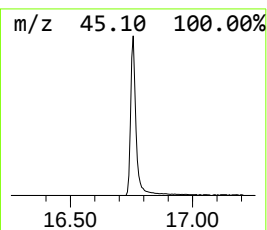
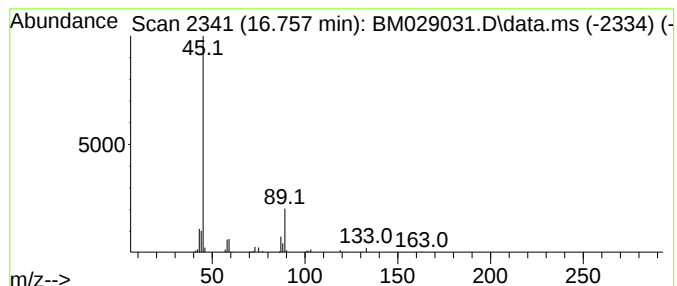
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Peak Number 5 Hexaethylene glycol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.756	19.30 ng	102130	Phenanthrene-d10	17.086

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexaethylene glycol	282	C12H26O7	002615-15-8	83
2			Tetraethylene glycol	194	C8H18O5	000112-60-7	83
3			Heptaethylene glycol	326	C14H30O8	005617-32-3	78
4			Ethanol, 2-[2-(ethenyloxy)ethoxy]-	132	C6H12O3	000929-37-3	78
5			Triethylene glycol	150	C6H14O4	000112-27-6	74



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VEOLIA-VNJ-208

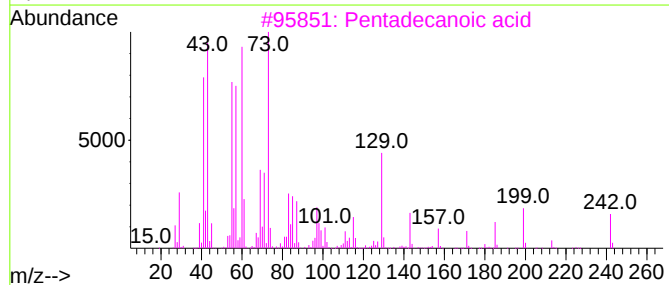
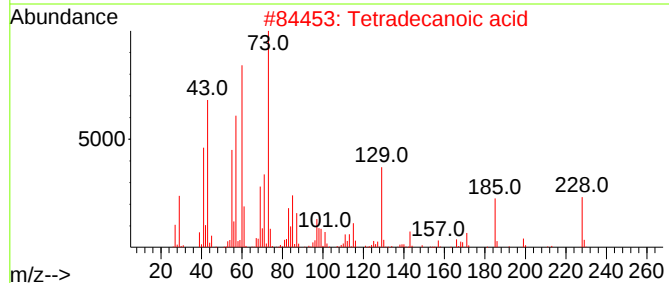
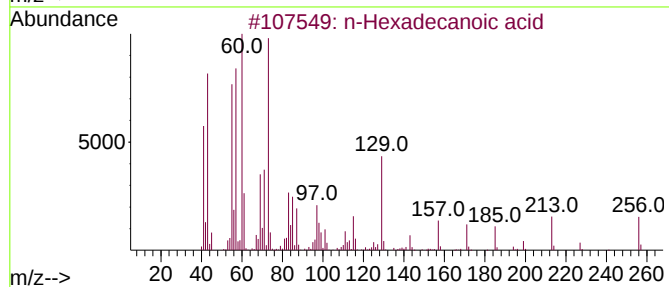
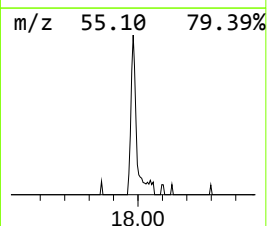
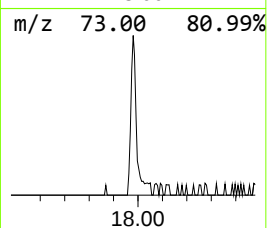
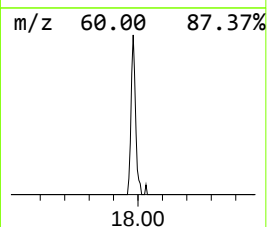
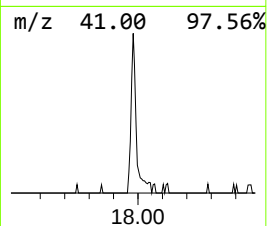
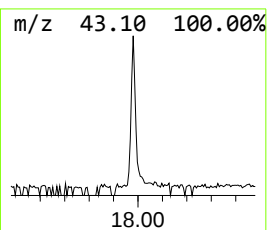
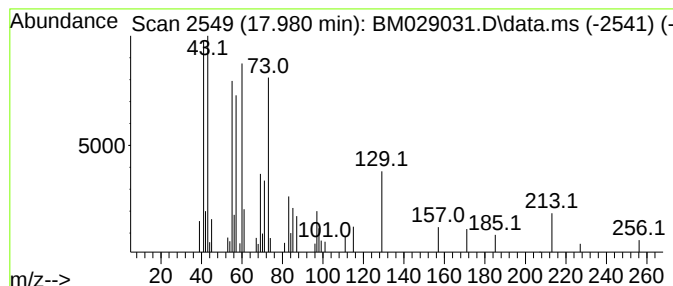
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.980	7.95 ng	42095	Phenanthrene-d10	17.086

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	72
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	64
4			Tridecanoic acid	214	C13H26O2	000638-53-9	53
5			Undecanoic acid	186	C11H22O2	000112-37-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
Data File : BM029031.D
Acq On : 08 Mar 2021 22:34
Operator : CG/JU
Sample : M1582-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
VEOLIA-VNJ-208

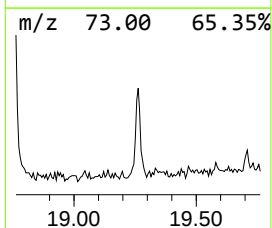
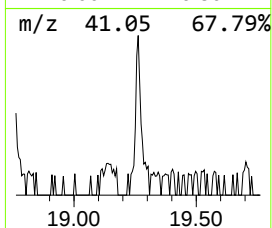
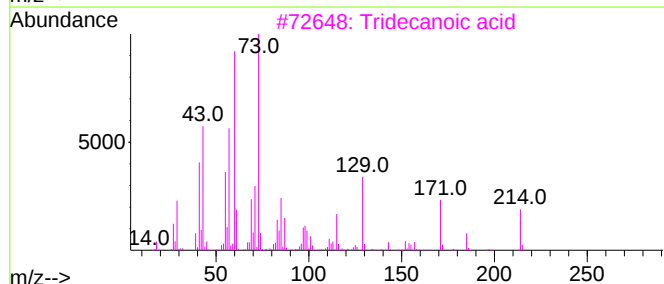
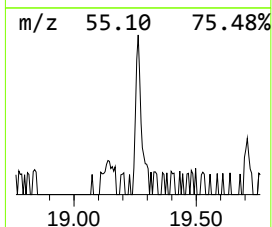
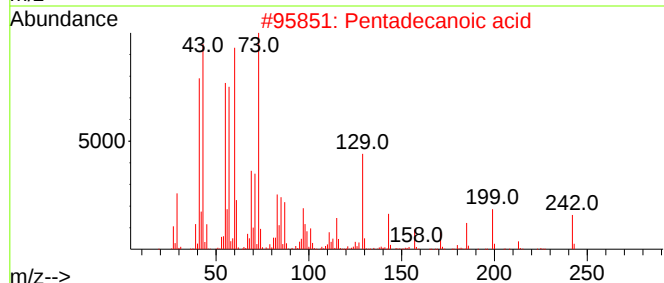
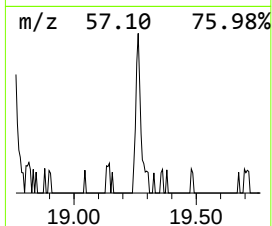
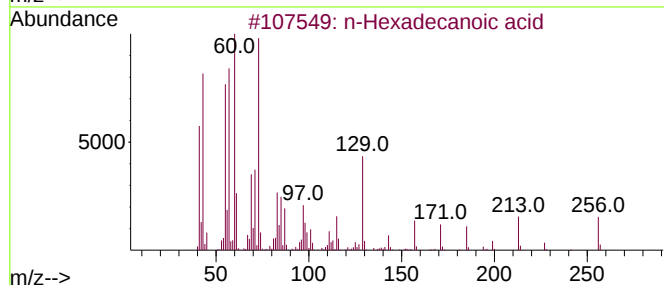
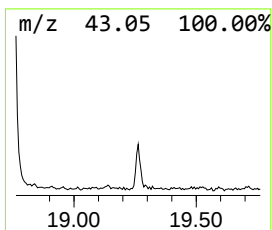
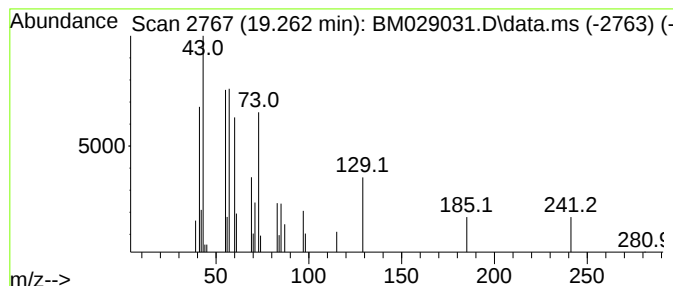
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pentadecanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.262	2.01 ng	15152	Chrysene-d12	21.256

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	72
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	72
3			Tridecanoic acid	214	C13H26O2	000638-53-9	53
4			Dodecanoic acid	200	C12H24O2	000143-07-7	50
5			n-Capric acid isopropyl ester	214	C13H26O2	002311-59-3	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
Data File : BM029031.D
Acq On : 08 Mar 2021 22:34
Operator : CG/JU
Sample : M1582-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
VEOLIA-VNJ-208

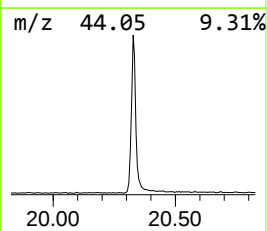
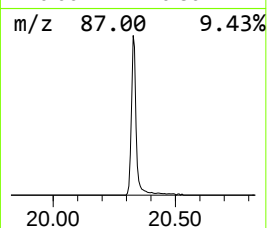
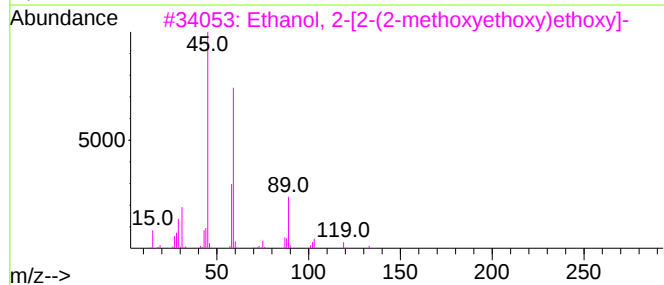
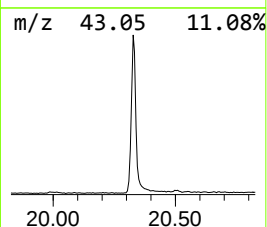
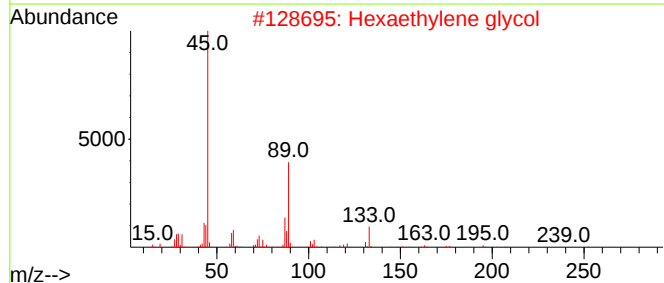
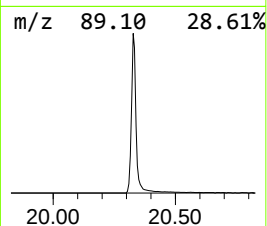
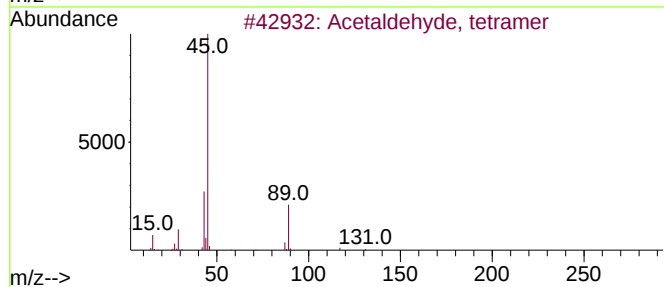
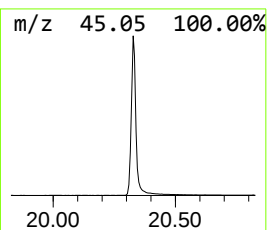
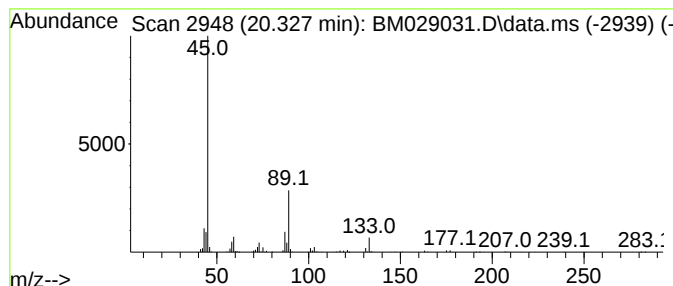
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Acetaldehyde, tetramer Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.327	91.84 ng	691713	Chrysene-d12	21.256

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetaldehyde, tetramer	176	C8H16O4	000108-62-3	72
2			Hexaethylene glycol	282	C12H26O7	002615-15-8	70
3			Ethanol, 2-[2-(2-methoxyethoxy)e...	164	C7H16O4	000112-35-6	64
4			Triethylene glycol	150	C6H14O4	000112-27-6	64
5			Heptaethylene glycol	326	C14H30O8	005617-32-3	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
Data File : BM029031.D
Acq On : 08 Mar 2021 22:34
Operator : CG/JU
Sample : M1582-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
VEOLIA-VNJ-208

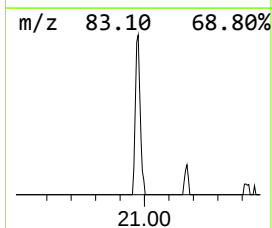
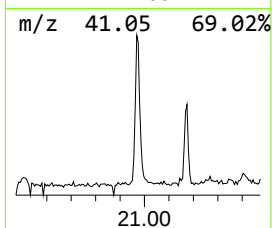
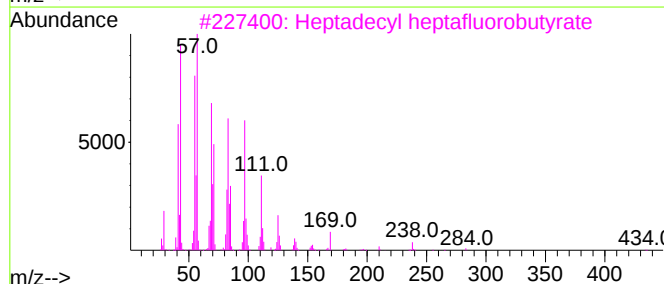
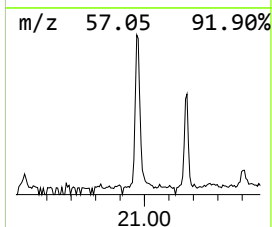
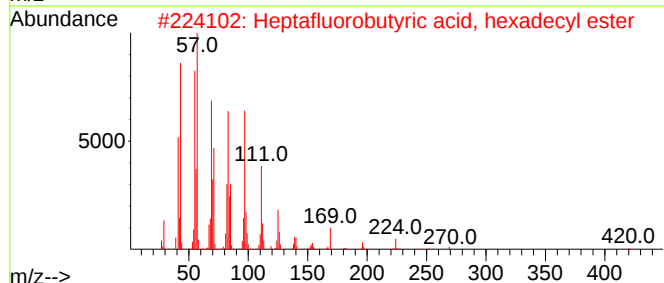
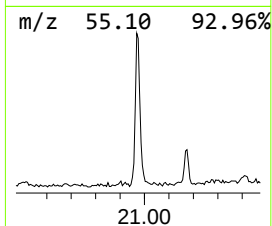
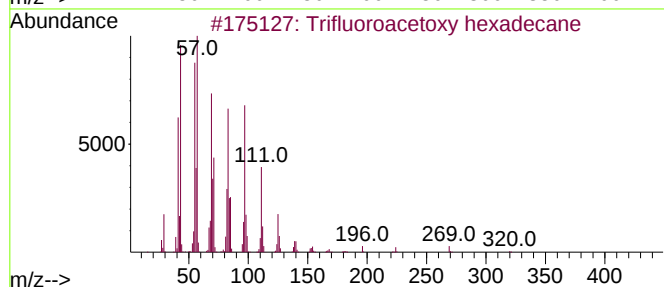
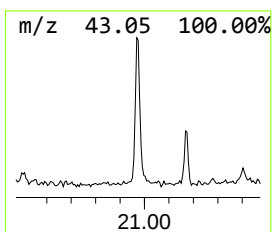
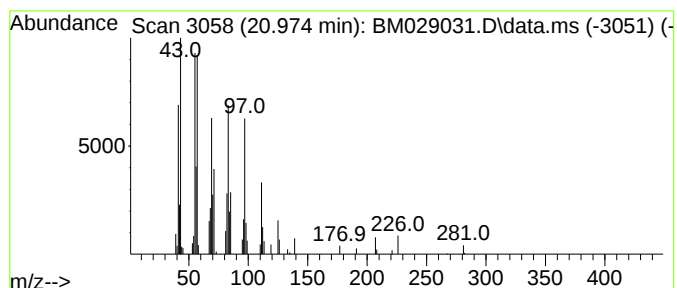
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Trifluoroacetoxy hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.974	8.06 ng	60703	Chrysene-d12	21.256

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	95
2			Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	95
3			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
4			1-Docosene	308	C22H44	001599-67-3	94
5			1-Heneicosanol	312	C21H44O	015594-90-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
Data File : BM029031.D
Acq On : 08 Mar 2021 22:34
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Sample : M1582-01
Misc :
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Instrument :
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ClientSampleId :
VEOLIA-VNJ-208

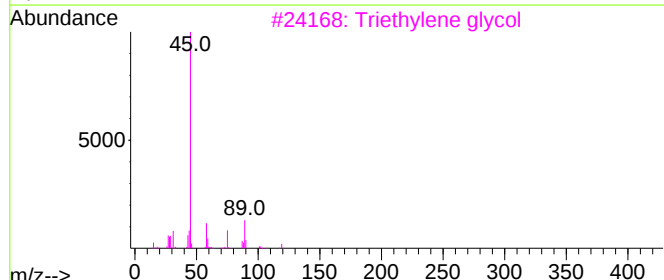
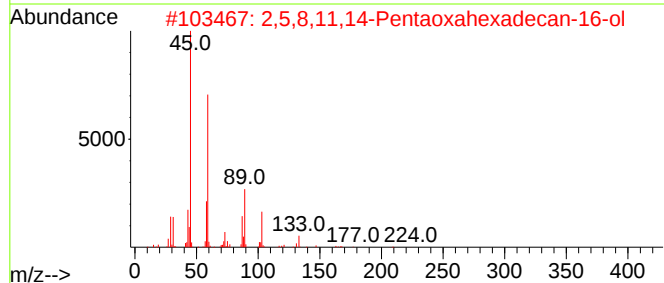
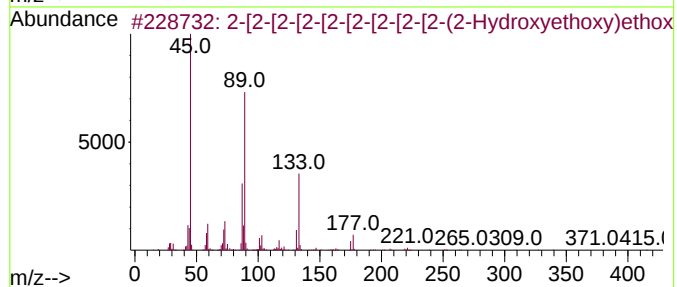
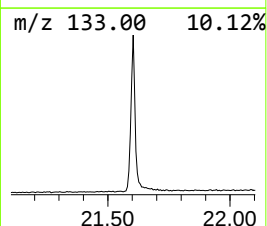
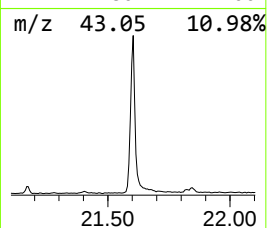
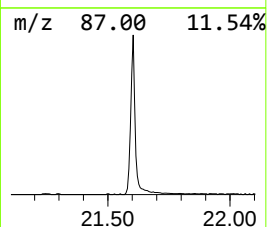
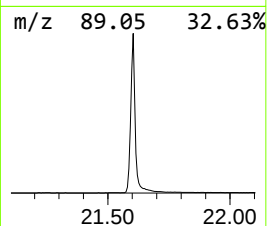
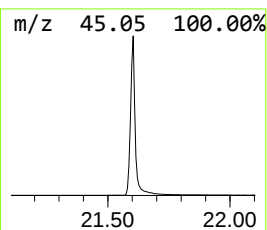
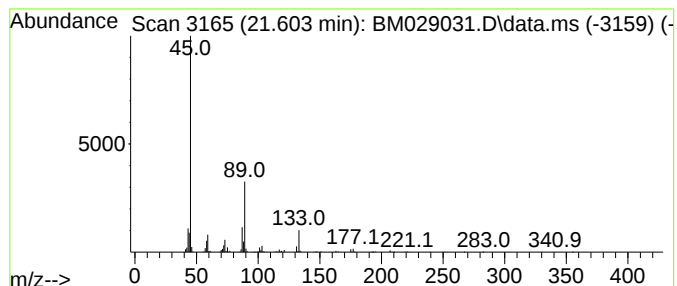
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2-[2-[2-[2-[2-[2-[2-(... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.603	136.70 ng	1029580	Chrysene-d12	21.256

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2	-[2-[2-[2-[2-[2-[2-(2-Hyd...	458	C20H42O11	1000352-12-6	90	
2	2,5,8,11,14-Pentaoxahexadecan-16-ol	252	C11H24O6	023778-52-1	78		
3	Triethylene glycol	150	C6H14O4	000112-27-6	72		
4	Heptaethylene glycol	326	C14H30O8	005617-32-3	70		
5	2-[2-[2-[2-[2-[2-(2-Hydroxyet...	370	C16H34O9	005117-19-1	62		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
Data File : BM029031.D
Acq On : 08 Mar 2021 22:34
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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
VEOLIA-VNJ-208

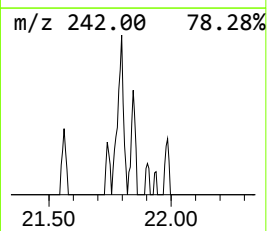
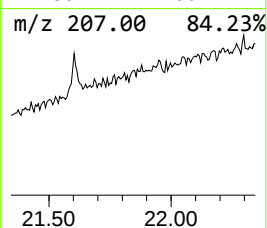
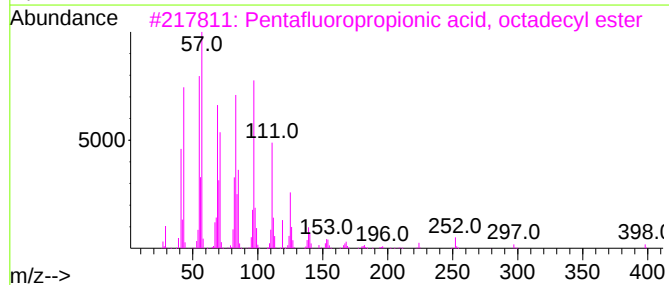
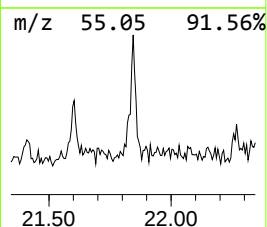
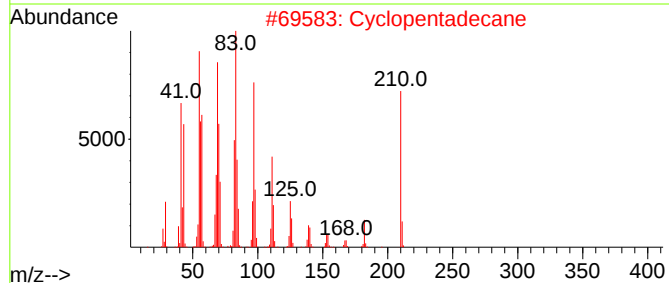
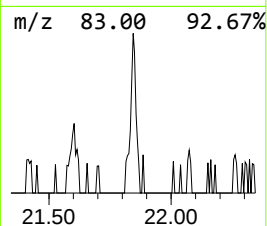
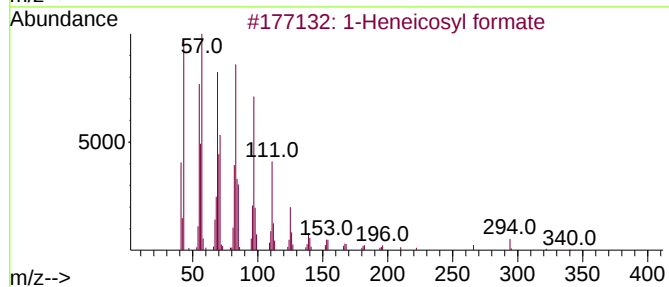
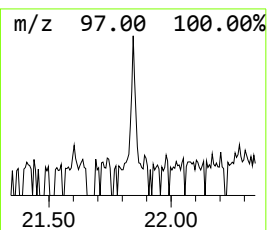
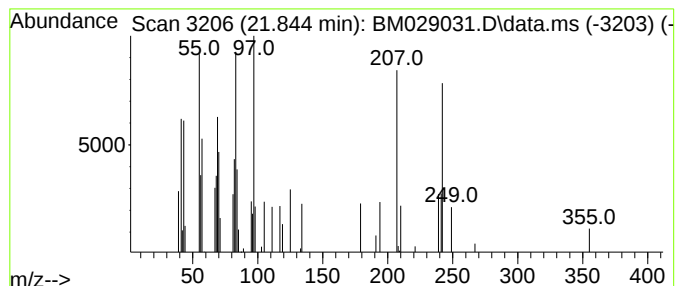
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown21.844 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.844	2.13 ng	16064	Chrysene-d12	21.256

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosyl formate	340	C22H44O2	077899-03-7	38
2			Cyclopentadecane	210	C15H30	000295-48-7	38
3			Pentafluoropropionic acid, octadecyl...	416	C21H37F5O2	959261-25-7	38
4			Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	35
5			Isoheptadecanol	256	C17H36O	057289-07-3	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
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Misc :
ALS Vial : 15 Sample Multiplier: 1

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ClientSampleId :
VEOLIA-VNJ-208

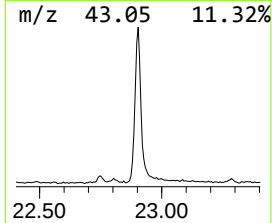
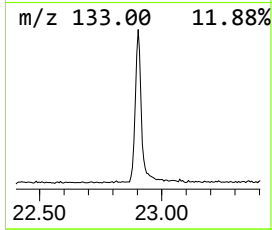
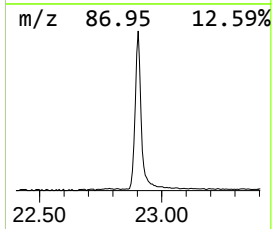
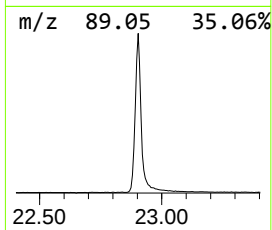
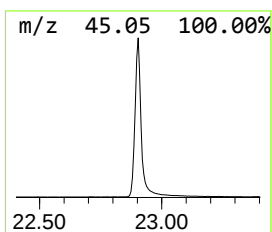
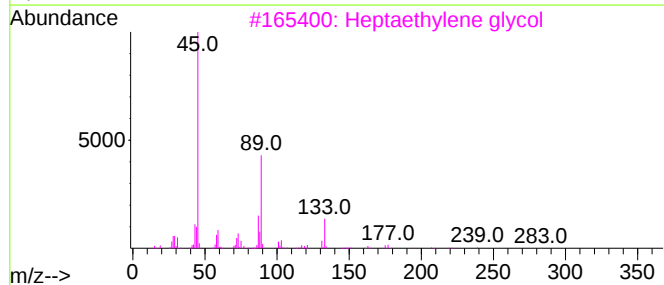
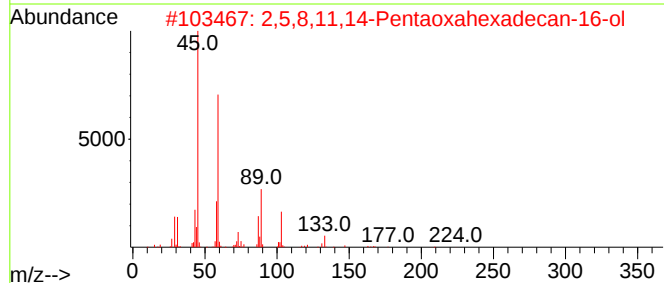
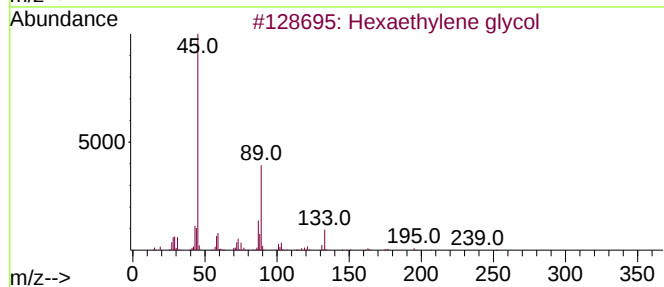
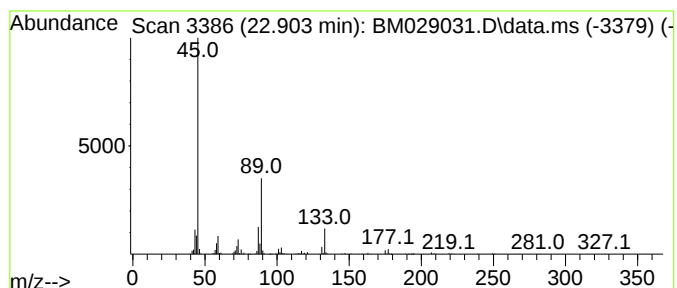
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 2,5,8,11,14-Pentaoxahexadec... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.903	140.74 ng	886112	Perylene-d12	23.415

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexaethylene glycol	282	C12H26O7	002615-15-8	83
2		2,5,8,11,14-Pentaoxahexadecan-16-ol	252	C11H24O6	023778-52-1	72
3		Heptaethylene glycol	326	C14H30O8	005617-32-3	70
4		2-[2-[2-[2-[2-[2-(2-Hydrox...	414	C18H38O10	1000352-12-5	70
5		2-[2-[2-[2-[2-[2-(2-Hydroxyet...	370	C16H34O9	005117-19-1	70



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Acq On : 08 Mar 2021 22:34
Operator : CG/JU
Sample : M1582-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
VEOLIA-VNJ-208

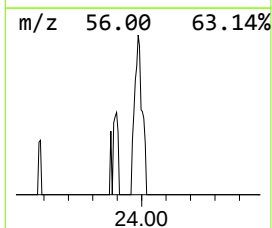
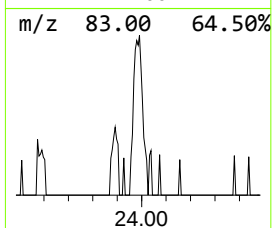
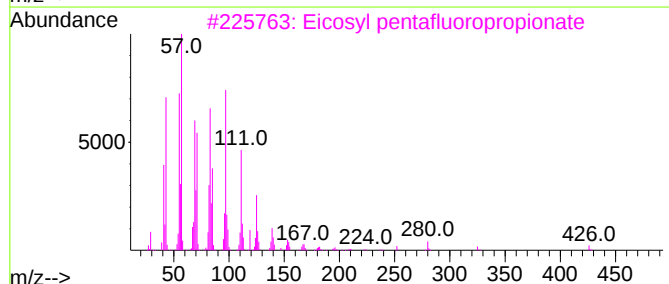
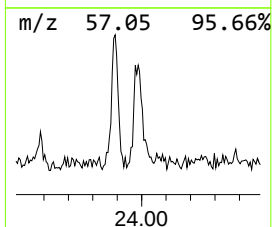
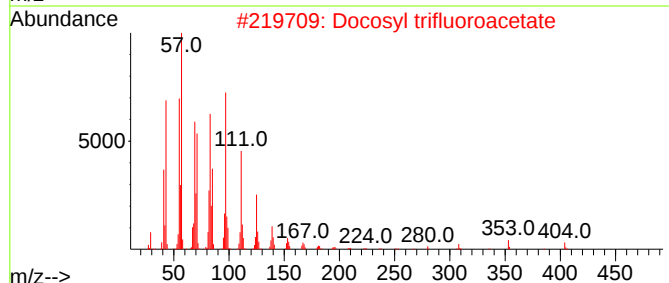
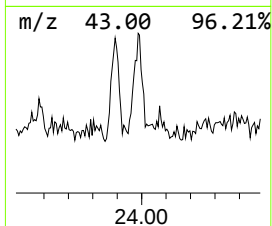
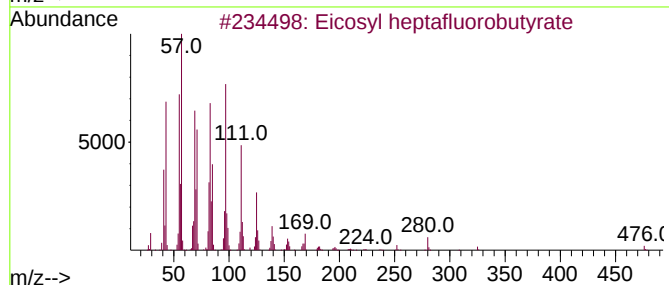
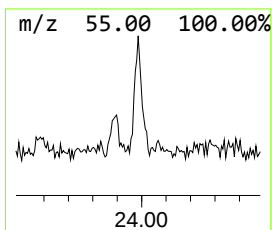
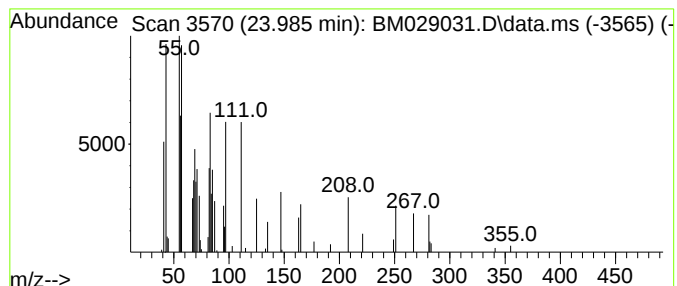
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown23.985 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.985	2.71 ng	17046	Perylene-d12	23.415

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosyl heptafluorobutyrate	494	C24H41F7O2	1000351-83-5	35
2			Docosyl trifluoroacetate	422	C24H45F3O2	1000351-75-5	35
3			Eicosyl pentafluoropropionate	444	C23H41F5O2	1000351-80-8	35
4			1-Hexacosanol	382	C26H54O	000506-52-5	30
5			n-Tetracosanol-1	354	C24H50O	000506-51-4	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
Data File : BM029031.D
Acq On : 08 Mar 2021 22:34
Operator : CG/JU
Sample : M1582-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

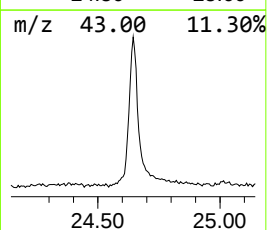
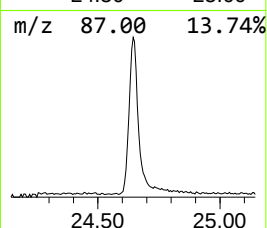
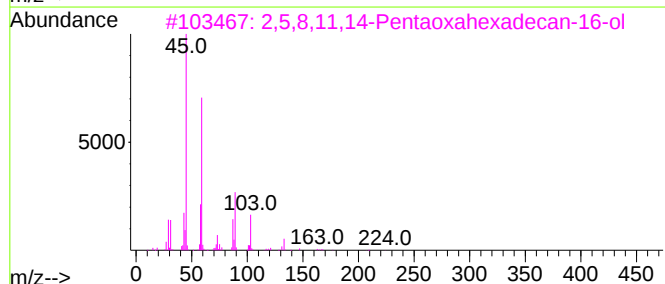
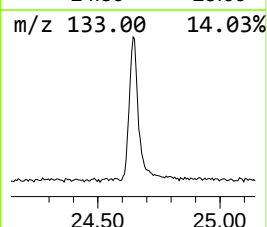
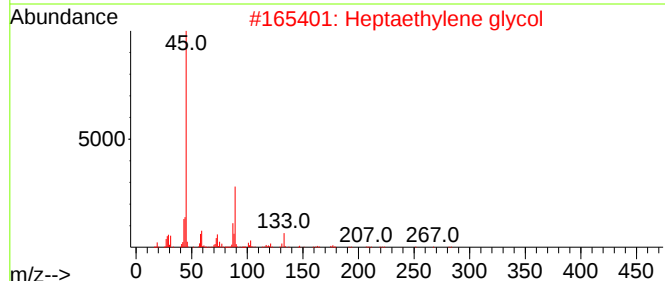
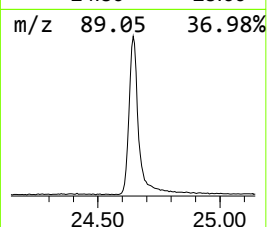
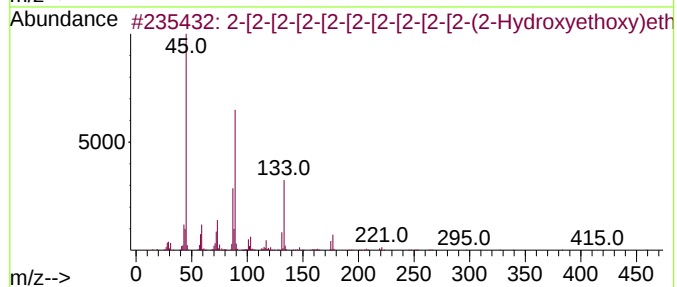
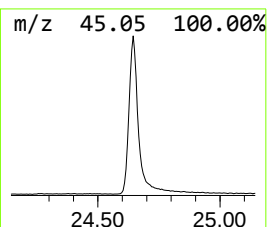
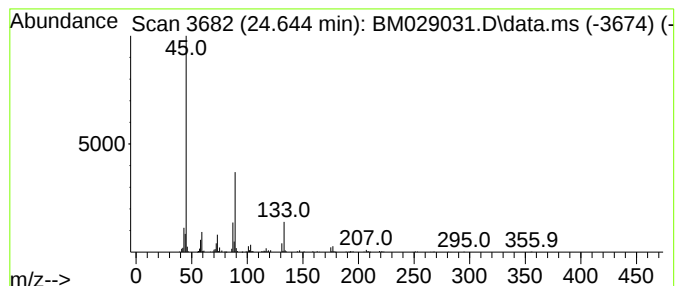
Instrument :
BNA_M
ClientSampleId :
VEOLIA-VNJ-208

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 15 2-[2-[2-[2-[2-[2-[2-[2-[... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.644	70.95 ng	446727	Perylene-d12	23.415	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-[2-[2-[2-[2-[2-[2-[2-(2-...	502	C22H46O12	1000352-13-2	91
2	Heptaethylene glycol	326	C14H30O8	005617-32-3	78
3	2,5,8,11,14-Pentaoxahexadecan-16-ol	252	C11H24O6	023778-52-1	72
4	Pentaethylene glycol	238	C10H22O6	004792-15-8	72
5	2-[2-[2-[2-[2-[2-[2-(2-Hydrox...	414	C18H38O10	1000352-12-5	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030821\
Data File : BM029031.D
Acq On : 08 Mar 2021 22:34
Operator : CG/JU
Sample : M1582-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
VEOLIA-VNJ-208

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
Propylene Glycol	3.399	241.0	ng	722818	1	7.669	59977	20.0
3-(2-Hydroxyeth...	13.310	33.6	ng	160526	3	14.339	95567	20.0
unknown14.227	14.227	2.4	ng	11361	3	14.339	95567	20.0
Tetraethylene g...	14.274	6.1	ng	29328	3	14.339	95567	20.0
Hexaethylene gl...	16.756	19.3	ng	102130	4	17.086	105836	20.0
n-Hexadecanoic ...	17.980	8.0	ng	42095	4	17.086	105836	20.0
2-[2-[2-[2-[2-[...	18.750	53.0	ng	280233	4	17.086	105836	20.0
Pentadecanoic acid	19.262	2.0	ng	15152	5	21.256	150638	20.0
Acetaldehyde, t...	20.327	91.8	ng	691713	5	21.256	150638	20.0
Trifluoroacetox...	20.974	8.1	ng	60703	5	21.256	150638	20.0
2-[2-[2-[2-[2-[...	21.603	136.7	ng	1029580	5	21.256	150638	20.0
unknown21.844	21.844	2.1	ng	16064	5	21.256	150638	20.0
2,5,8,11,14-Pen...	22.903	140.7	ng	886112	6	23.415	125926	20.0
unknown23.985	23.985	2.7	ng	17046	6	23.415	125926	20.0
2-[2-[2-[2-[2-[...	24.644	71.0	ng	446727	6	23.415	125926	20.0