

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082521\
 Data File : BG049969.D
 Acq On : 25 Aug 2021 15:31
 Operator : CG/JU
 Sample : M3506-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 N29305

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-BG080321.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG049969.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.534	419	425	433	rVB	180178	301397	17.62%	3.868%
2	7.038	676	681	689	rVB2	11006	20467	1.20%	0.263%
3	7.156	694	701	709	rBV	114735	206862	12.09%	2.655%
4	7.978	834	841	848	rVV	95941	162068	9.47%	2.080%
5	8.066	848	856	863	rVB5	9029	19690	1.15%	0.253%
6	9.147	1032	1040	1049	rBV	331196	591802	34.59%	7.595%
7	9.447	1086	1091	1101	rVB4	18700	30305	1.77%	0.389%
8	9.852	1154	1160	1171	rBV4	27554	69978	4.09%	0.898%
9	10.557	1274	1280	1289	rBV2	81369	145192	8.49%	1.863%
10	10.798	1314	1321	1329	rBV	126898	242637	14.18%	3.114%
11	11.409	1420	1425	1431	rBV2	20520	39305	2.30%	0.504%
12	11.702	1469	1475	1481	rBV	31844	54428	3.18%	0.699%
13	11.761	1481	1485	1491	rVB2	28002	49268	2.88%	0.632%
14	11.943	1510	1516	1522	rBV2	21559	36431	2.13%	0.468%
15	13.253	1731	1739	1746	rBV	868277	1346697	78.72%	17.284%
16	14.275	1908	1913	1918	rBV7	10050	22360	1.31%	0.287%
17	14.634	1968	1974	1981	rBV	202366	318087	18.59%	4.082%
18	14.734	1987	1991	1996	rBV2	18772	27531	1.61%	0.353%
19	15.045	2040	2044	2050	rBV6	22113	35641	2.08%	0.457%
20	15.450	2108	2113	2119	rBV9	11810	23147	1.35%	0.297%
21	16.132	2223	2229	2238	rVB	740573	1148352	67.12%	14.738%
22	17.389	2438	2443	2450	rVB	230495	356978	20.87%	4.582%
23	17.700	2492	2496	2504	rVB9	16944	31900	1.86%	0.409%
24	18.041	2551	2554	2559	rBV5	29414	41350	2.42%	0.531%
25	18.282	2591	2595	2598	rBV6	24190	33306	1.95%	0.427%
26	20.003	2883	2888	2894	rVB	1406920	1710774	100.00%	21.957%
27	21.671	3166	3172	3176	rVB2	205595	321613	18.80%	4.128%
28	24.908	3716	3723	3736	rVB3	134015	404073	23.62%	5.186%

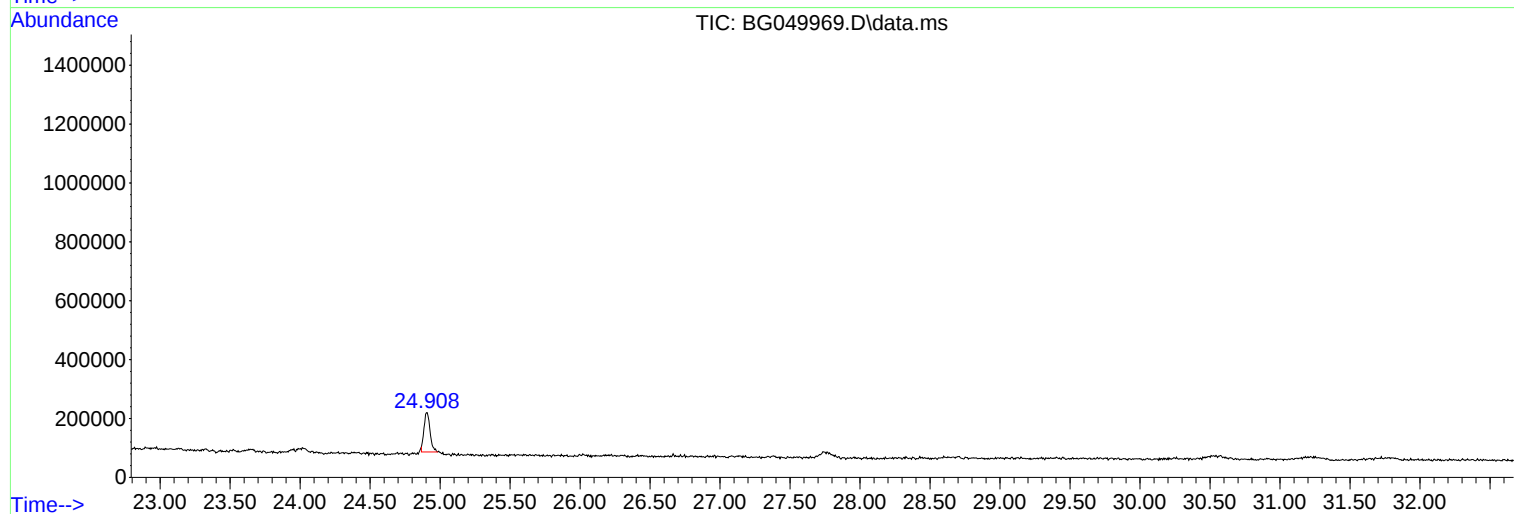
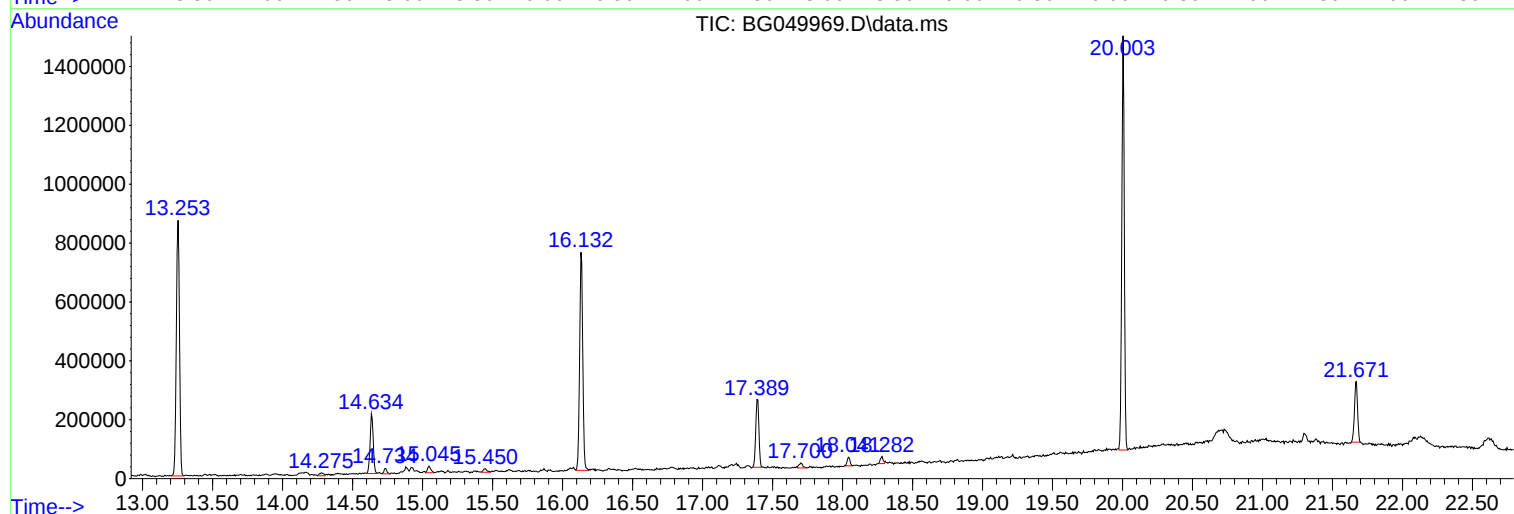
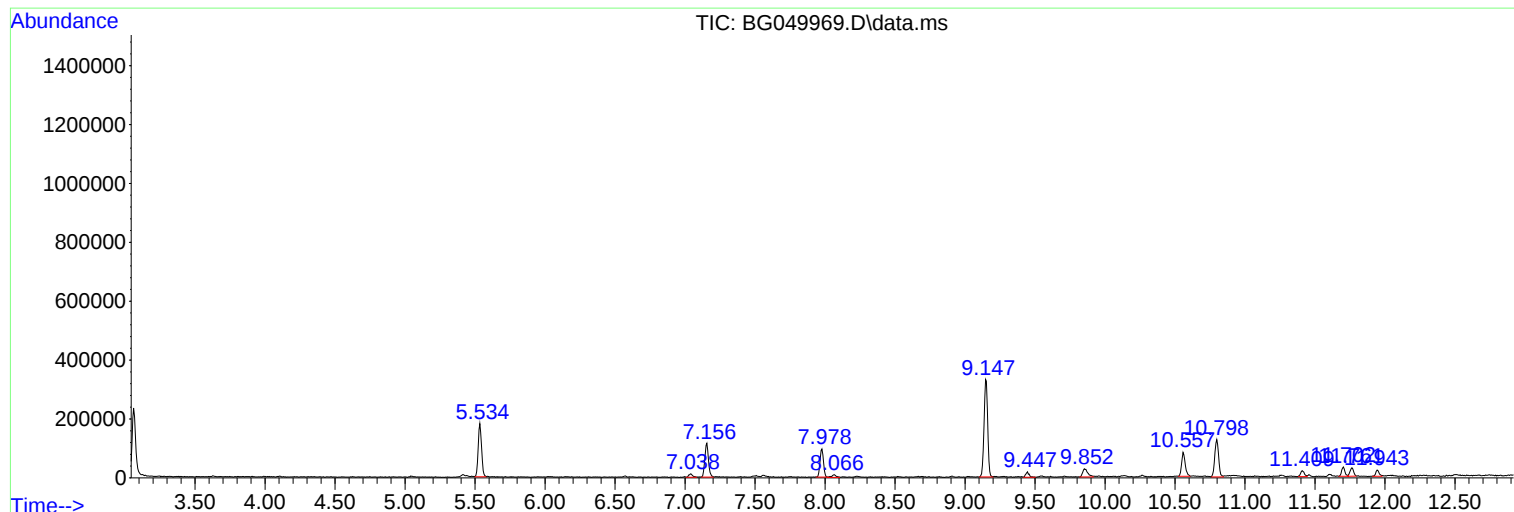
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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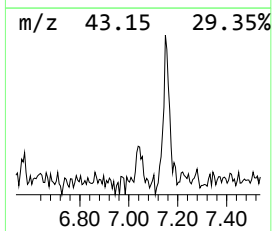
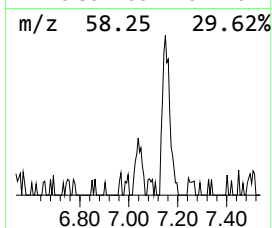
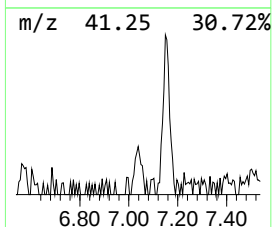
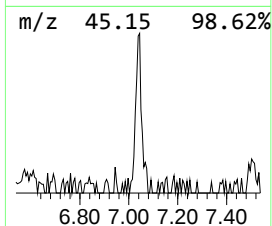
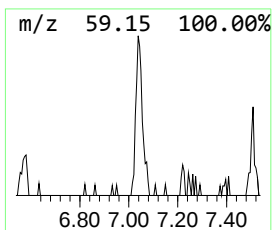
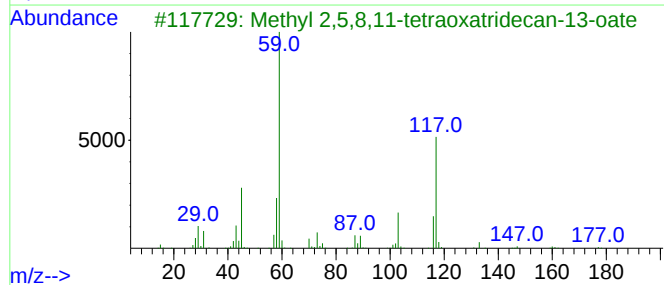
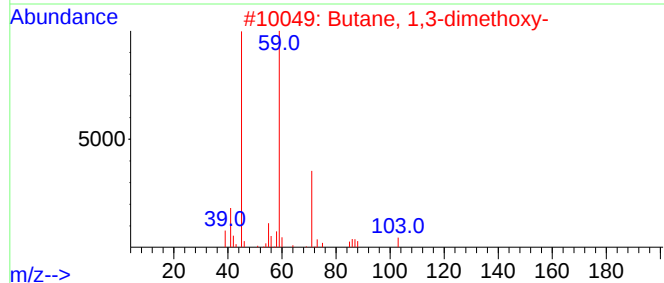
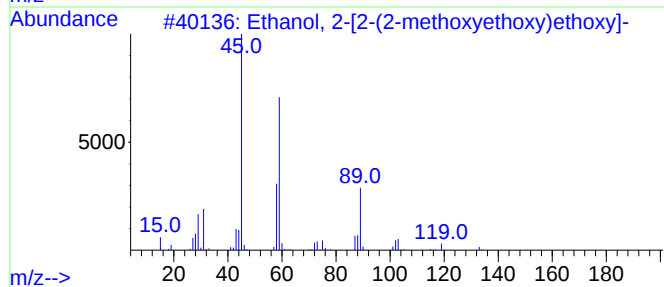
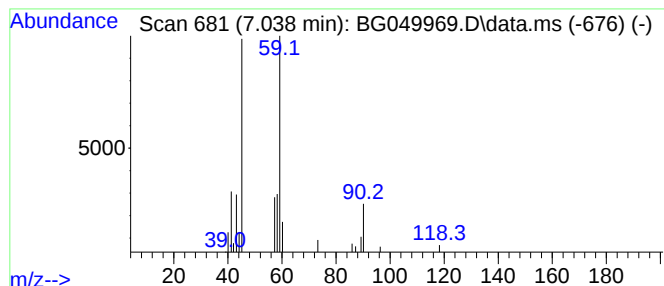
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethanol, 2-[2-(2-methoxyeth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.038	2.53 ng	20467	1,4-Dichlorobenzene-d4	7.978

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-[2-(2-methoxyethoxy)e...	164	C7H16O4	000112-35-6	64
2			Butane, 1,3-dimethoxy-	118	C6H14O2	010143-66-5	36
3			Methyl 2,5,8,11-tetraoxatridecan...	236	C10H20O6	1000366-78-2	28
4			3-Methoxypropionic acid	104	C4H8O3	002544-06-1	23
5			2-Hydroxybutanoic acid hydrazide	118	C4H10N2O2	1010197-30-7	9



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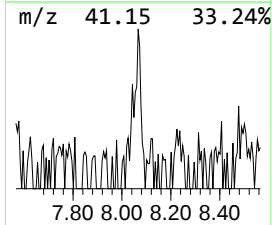
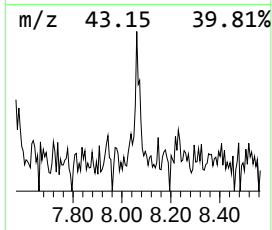
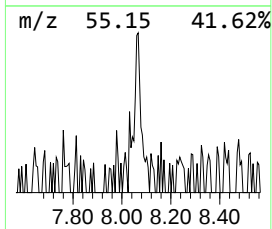
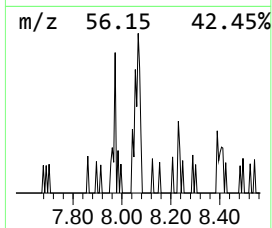
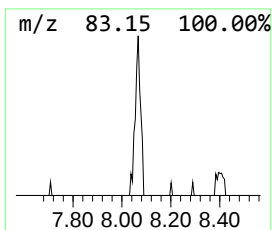
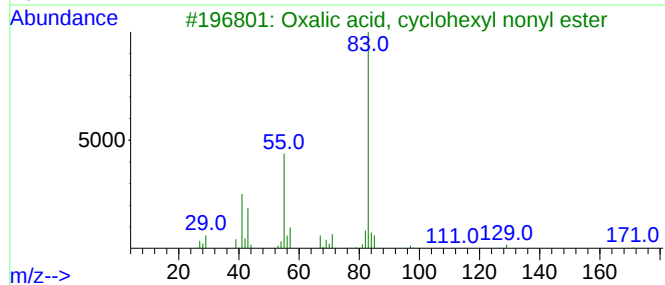
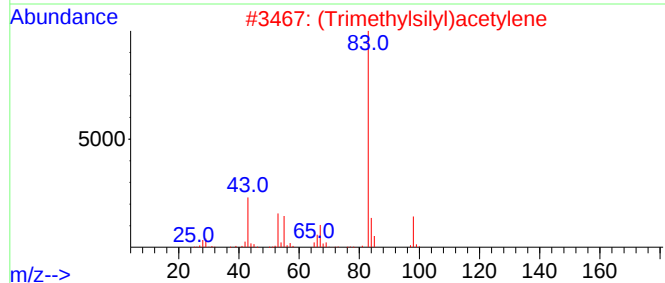
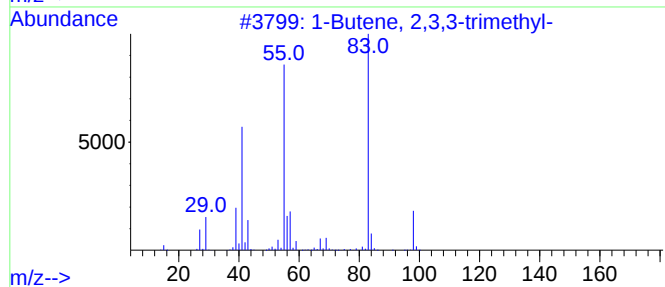
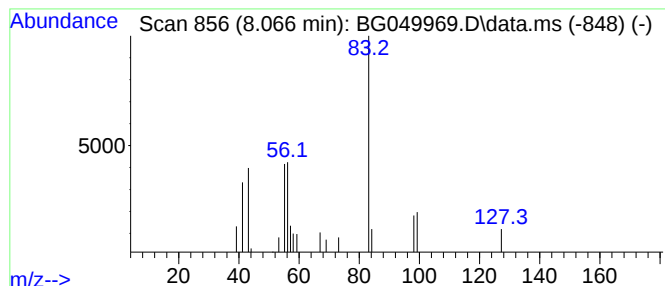
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 unknown8.066 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.066	2.43 ng	19690	1,4-Dichlorobenzene-d4	7.978

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Butene, 2,3,3-trimethyl-	98	C7H14	000594-56-9	47
2			(Trimethylsilyl)acetylene	98	C5H10Si	001066-54-2	43
3			Oxalic acid, cyclohexyl nonyl ester	298	C17H30O4	1000309-31-1	38
4			Oxalic acid, cyclohexyl dodecyl ...	340	C20H36O4	1000309-31-4	38
5			Propanenitrile, 2-(dimethylamino)-	98	C5H10N2	005350-67-4	36



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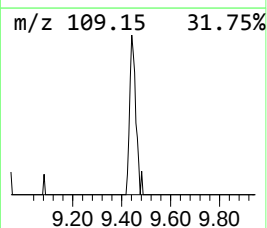
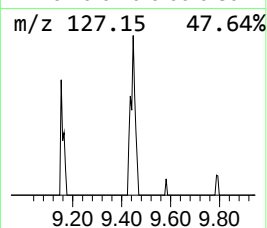
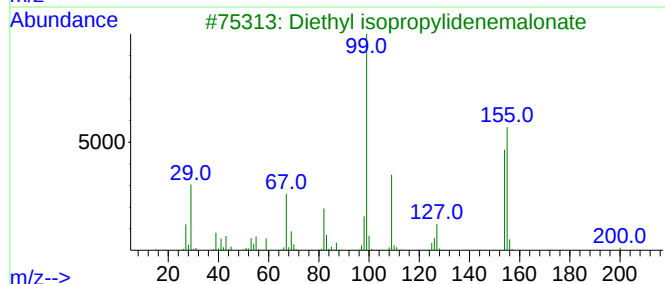
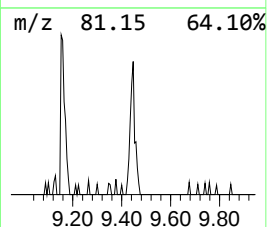
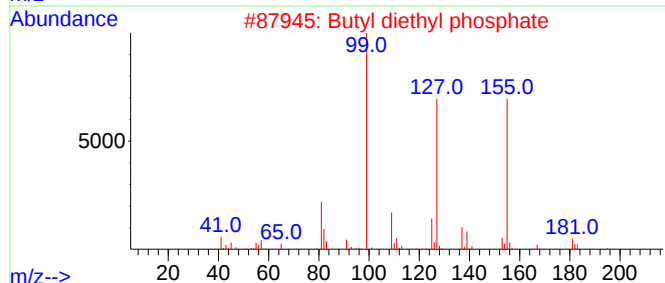
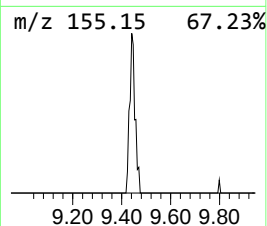
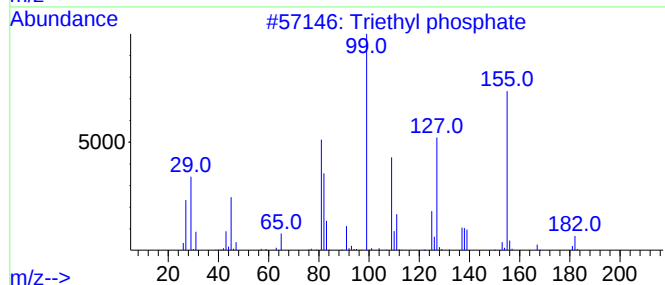
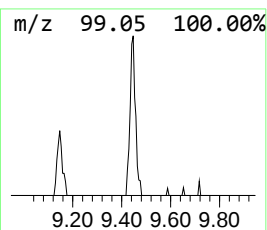
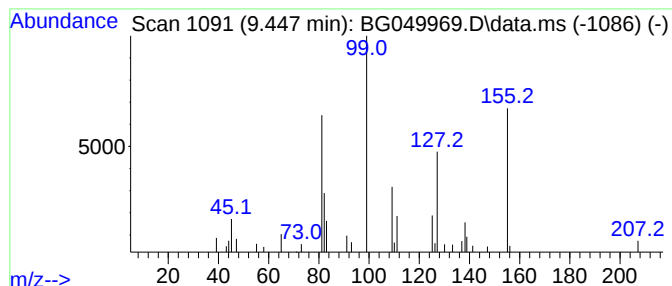
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Triethyl phosphate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.447	2.50 ng	30305	Naphthalene-d8	10.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triethyl phosphate	182	C6H15O4P	000078-40-0	90
2			Butyl diethyl phosphate	210	C8H19O4P	1000298-58-0	56
3			Diethyl isopropylidenemalonate	200	C10H16O4	006802-75-1	42
4			Phosphoric acid, monoethyl ester	126	C2H7O4P	001623-14-9	35
5			Diethyl 1,1-cyclobutanedicarboxy...	200	C10H16O4	003779-29-1	33



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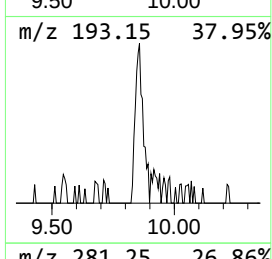
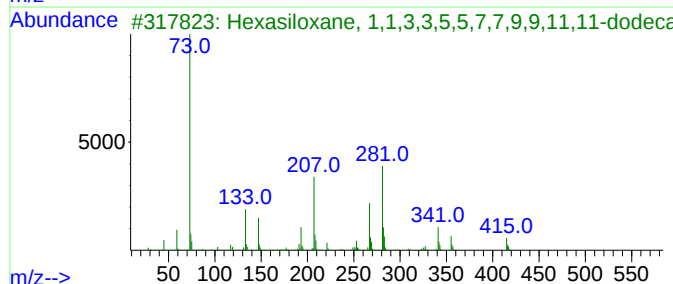
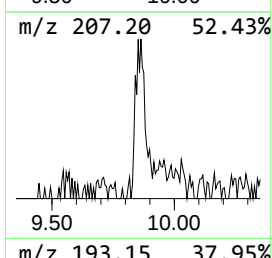
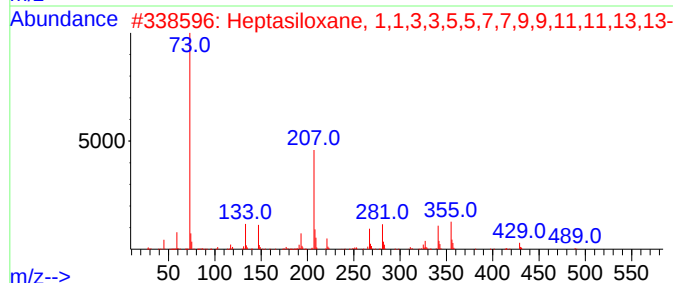
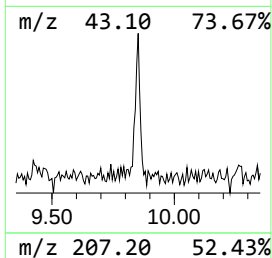
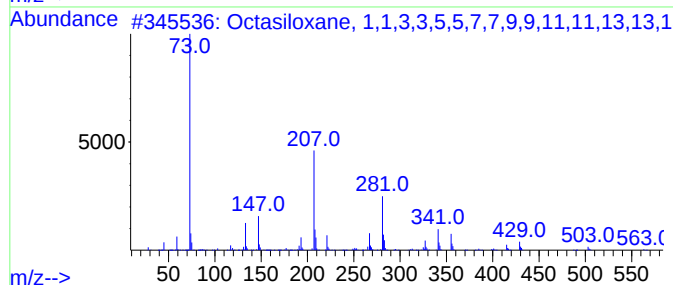
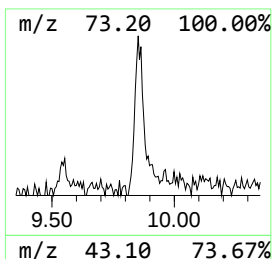
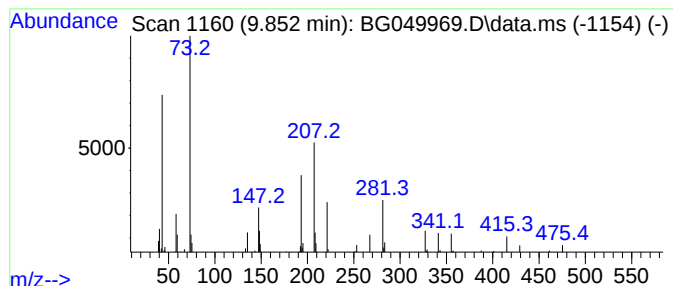
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Peak Number 4 Octasiloxane, 1,1,3,3,5,5,7... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.852	5.77 ng	69978	Naphthalene-d8	10.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octasiloxane, 1,1,3,3,5,5,7,7,9,...	578	C16H5007Si8	019095-24-0	64
2			Heptasiloxane, 1,1,3,3,5,5,7,7,9...	504	C14H4406Si7	019095-23-9	49
3			Hexasiloxane, 1,1,3,3,5,5,7,7,9...	430	C12H3805Si6	000995-82-4	40
4			1H-Indole-2-carboxylic acid, 6-(...	355	C21H25NO4	1010316-17-5	27
5			Silicic acid, diethyl bis(trimet...	296	C10H2804Si3	003555-45-1	27



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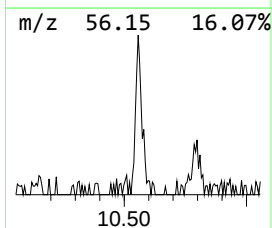
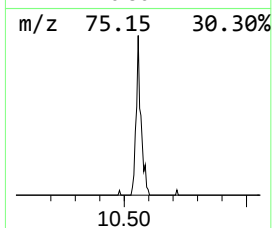
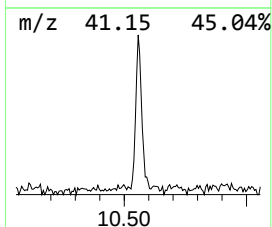
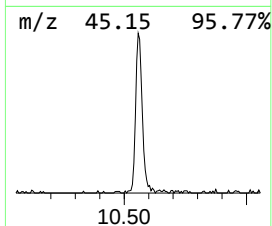
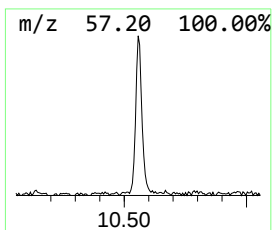
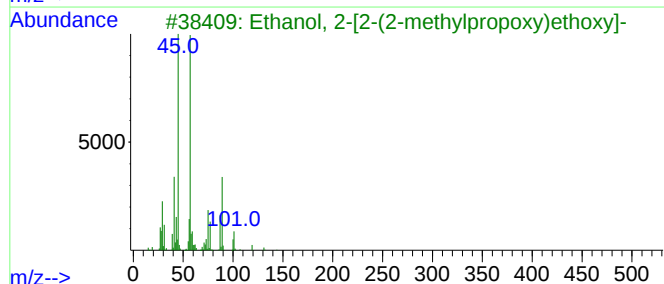
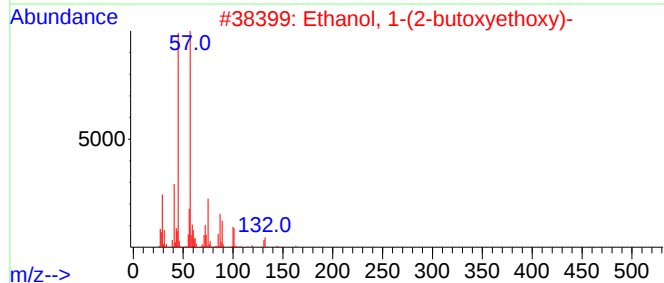
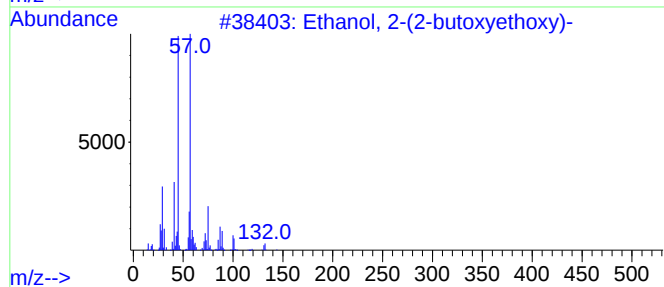
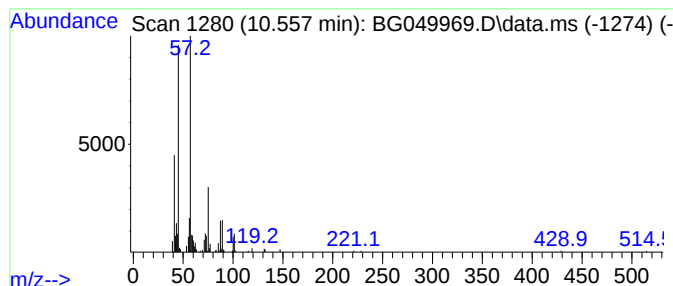
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Peak Number 5 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.557	11.97 ng	145192	Naphthalene-d8	10.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	78
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	78
3			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	72
4			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
5			Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	50



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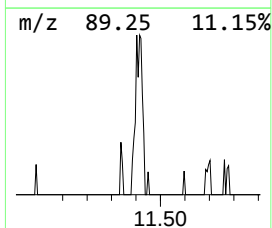
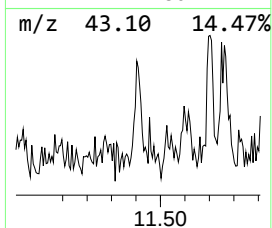
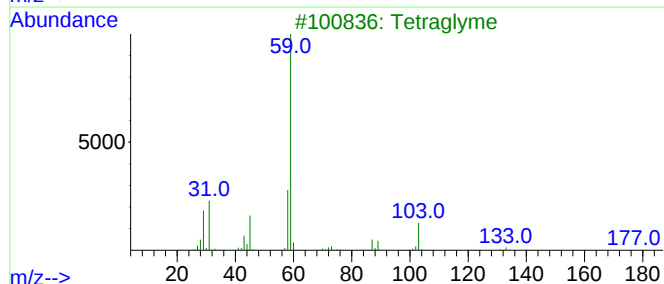
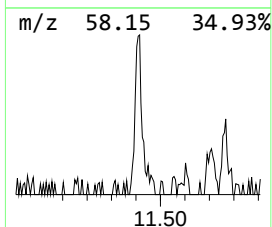
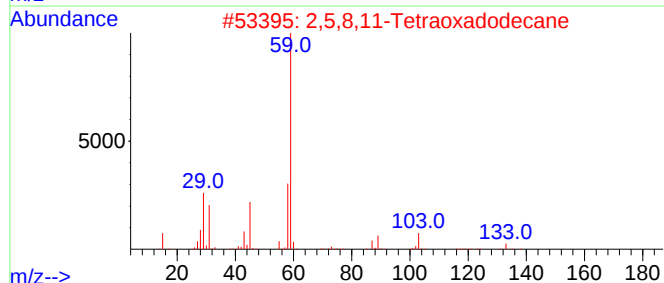
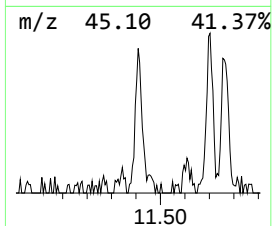
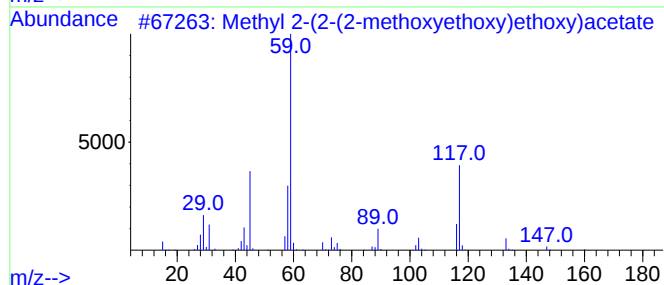
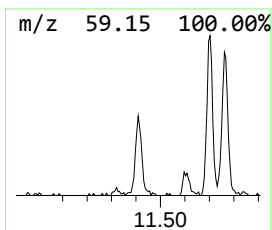
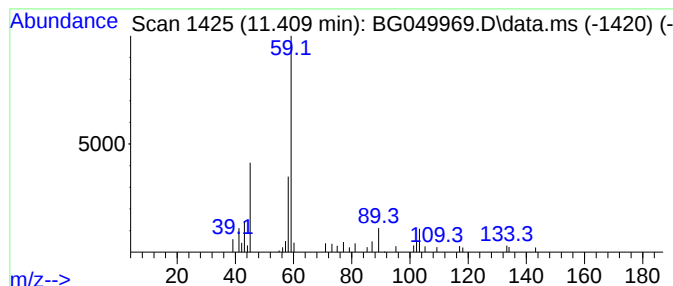
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Methyl 2-(2-(2-methoxyethoxy)ethoxy)acetate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.409	3.24 ng	39305	Naphthalene-d8	10.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 2-(2-(2-methoxyethoxy)ethoxy)acetate	192	C8H16O5	1000366-88-2	64
2			2,5,8,11-Tetraoxadodecane	178	C8H18O4	000112-49-2	64
3			Tetraglyme	222	C10H22O5	000143-24-8	59
4			Ethane, 1,1'-oxybis[2-methoxy-	134	C6H14O3	000111-96-6	56
5			Isopropyldimethylsilane	102	C5H14Si	1000433-69-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082521\
Data File : BG049969.D
Acq On : 25 Aug 2021 15:31
Operator : CG/JU
Sample : M3506-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
N29305

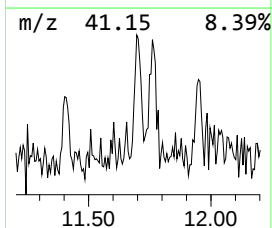
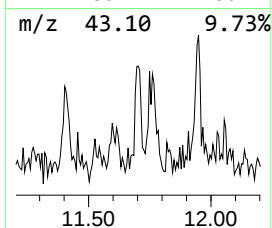
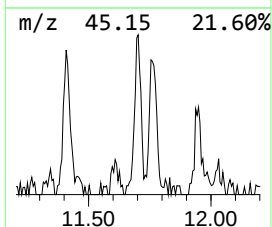
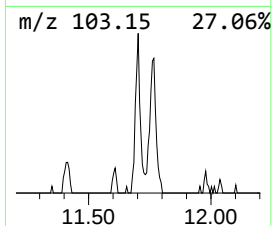
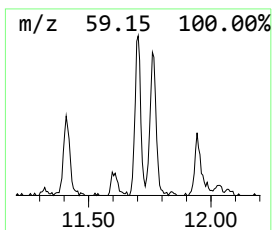
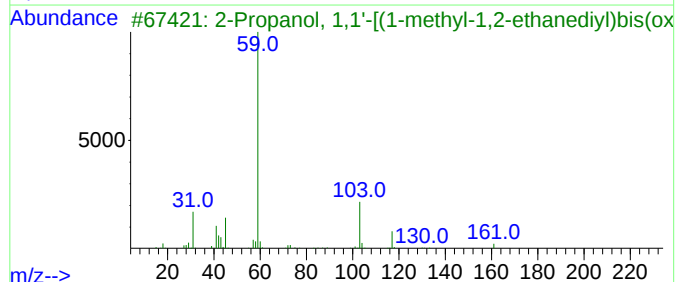
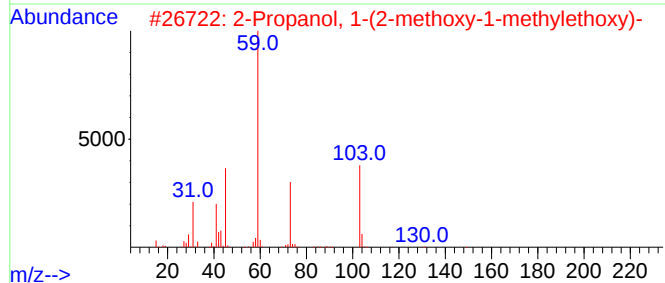
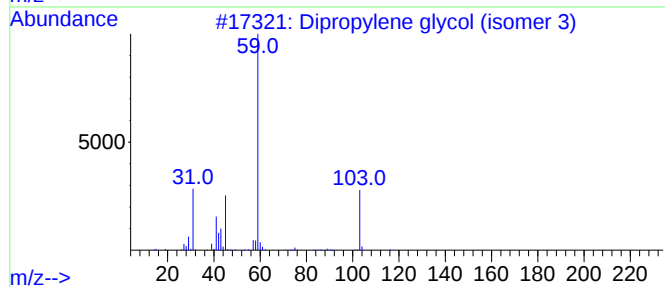
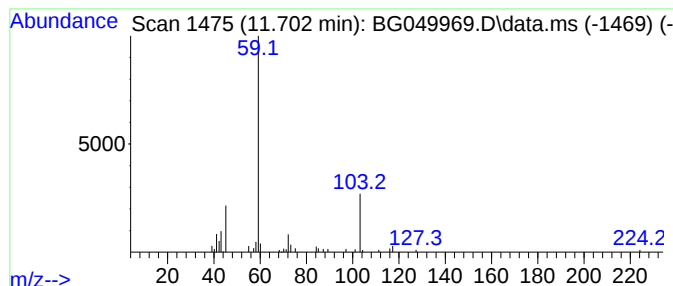
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Dipropylene glycol (isomer 3) Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.702	4.49 ng	54428	Naphthalene-d8	10.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dipropylene glycol (isomer 3)	134	C6H14O3	1000506-25-5	72
2			2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	72
3			2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	64
4			Tetraglyme	222	C10H22O5	000143-24-8	64
5			2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082521\
Data File : BG049969.D
Acq On : 25 Aug 2021 15:31
Operator : CG/JU
Sample : M3506-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
N29305

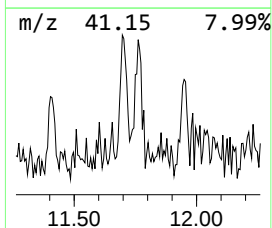
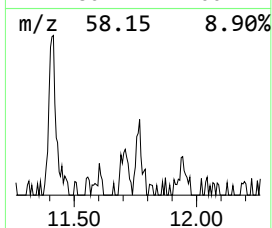
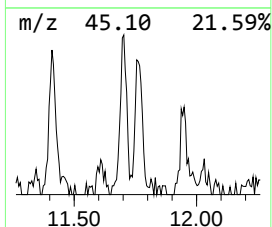
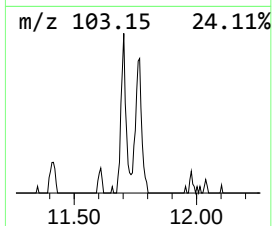
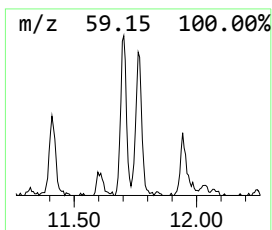
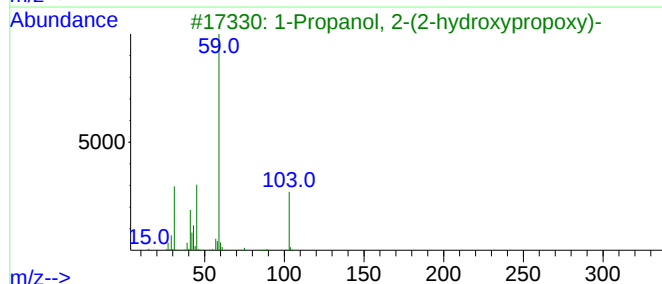
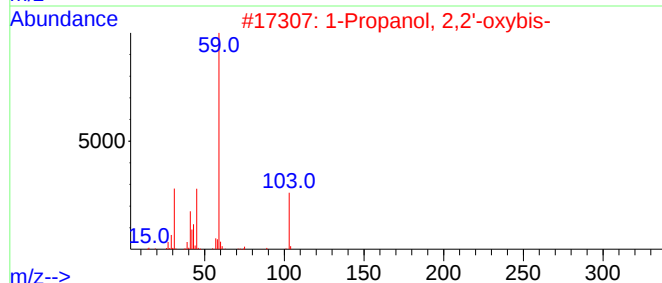
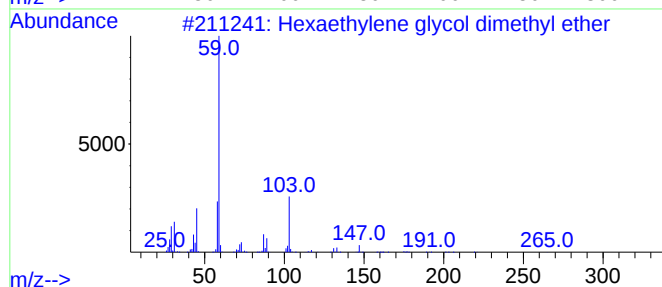
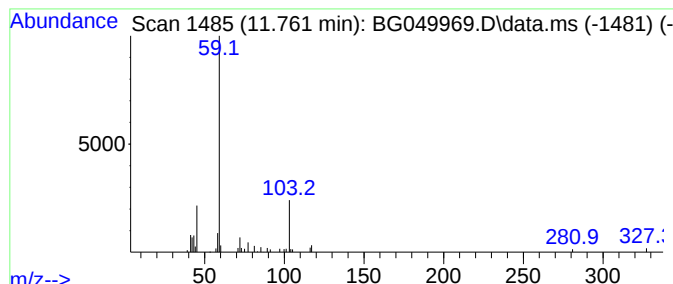
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Hexaethylene glycol dimethy... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.761	4.06 ng	49268	Naphthalene-d8	10.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexaethylene glycol dimethyl ether	310	C14H30O7	001072-40-8	78
2			1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	72
3			1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72
4			1-(1-tert-Butoxypropan-2-yloxy)p...	190	C10H22O3	132739-31-2	64
5			Propane, 1,2-dimethoxy-	104	C5H12O2	007778-85-0	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082521\
Data File : BG049969.D
Acq On : 25 Aug 2021 15:31
Operator : CG/JU
Sample : M3506-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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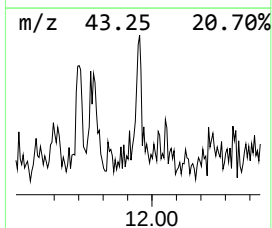
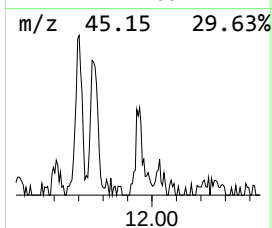
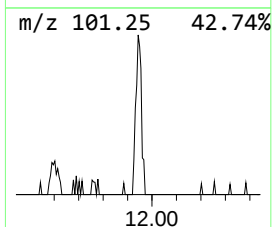
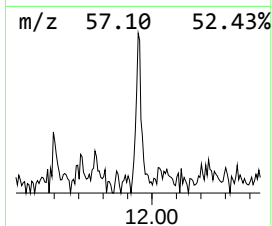
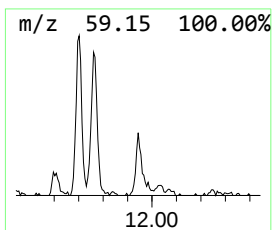
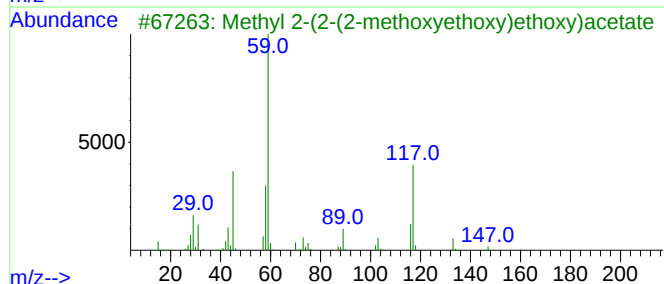
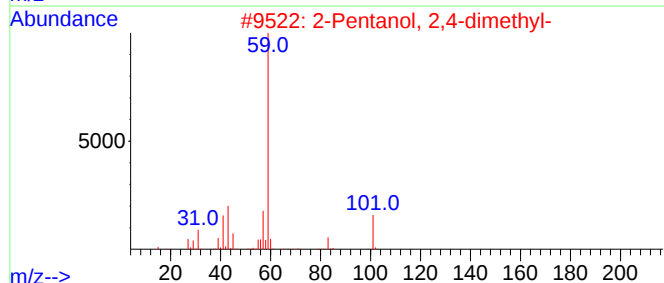
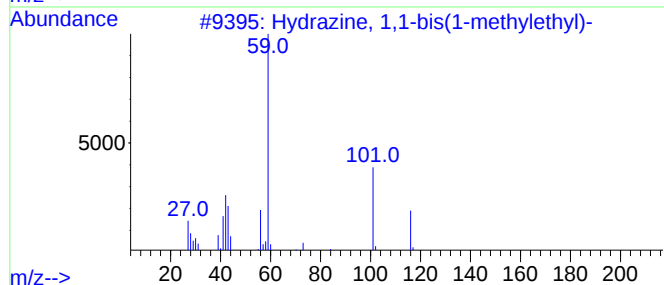
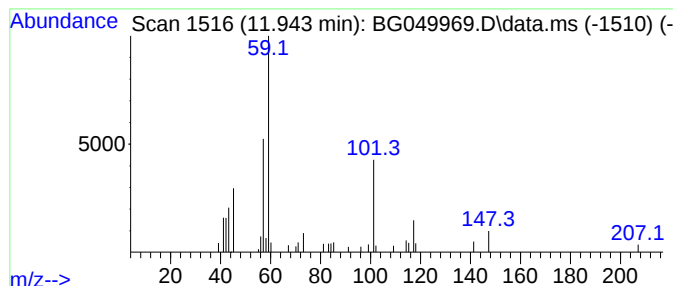
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Hydrazine, 1,1-bis(1-methyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.943	3.00 ng	36431	Naphthalene-d8	10.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	50
2			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	40
3			Methyl 2-(2-(2-methoxyethoxy)ethoxy)acetate	192	C8H16O5	1000366-88-2	38
4			Methyl 2,5,8,11,14,17,20,23-octa...	412	C18H36O10	1000366-81-2	38
5			2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082521\
Data File : BG049969.D
Acq On : 25 Aug 2021 15:31
Operator : CG/JU
Sample : M3506-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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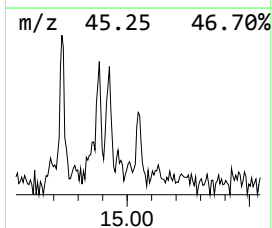
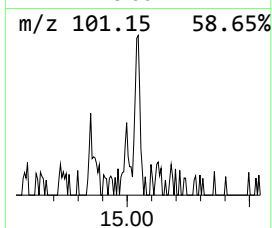
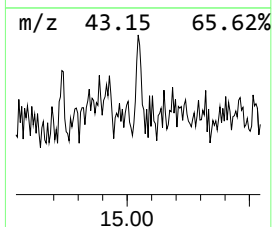
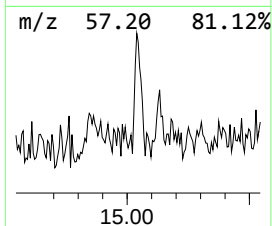
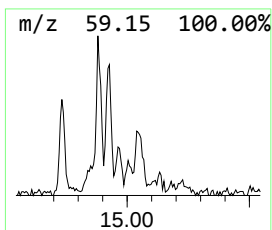
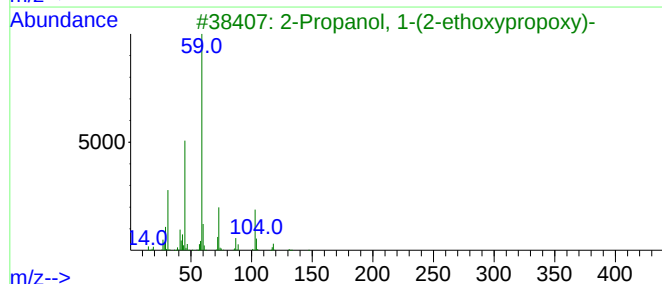
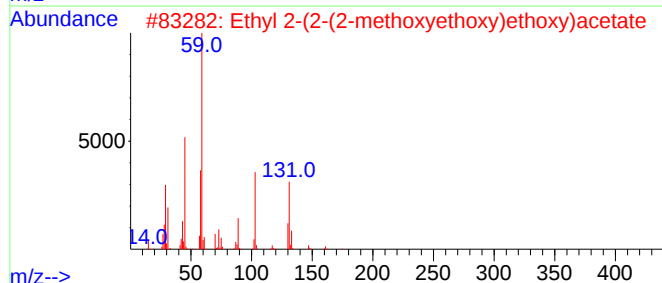
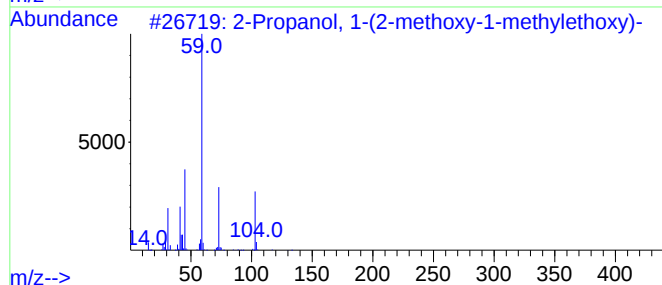
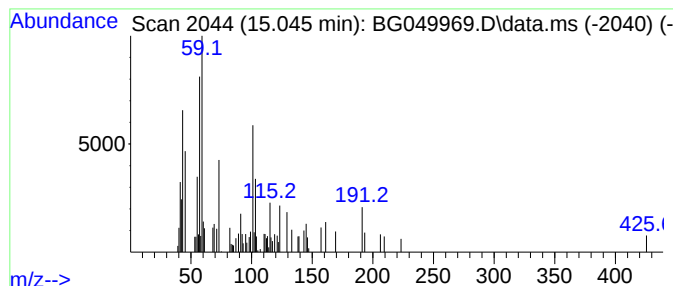
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown15.045 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.045	2.24 ng	35641	Acenaphthene-d10	14.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	43
2			Ethyl 2-(2-(2-methoxyethoxy)etho...	206	C9H18O5	1000366-89-0	43
3			2-Propanol, 1-(2-ethoxypropoxy)-	162	C8H18O3	010143-32-5	37
4			Propanedioic acid, dimethyl ester	132	C5H8O4	000108-59-8	32
5			n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082521\
Data File : BG049969.D
Acq On : 25 Aug 2021 15:31
Operator : CG/JU
Sample : M3506-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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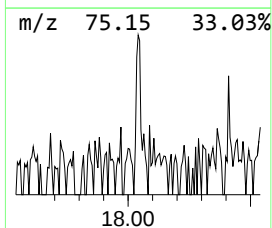
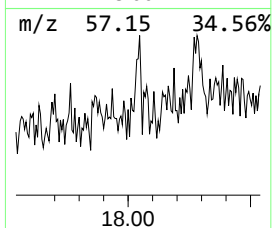
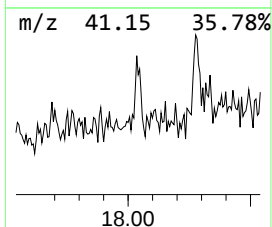
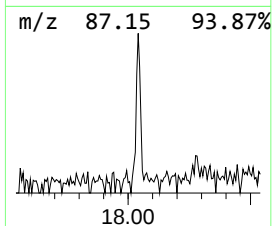
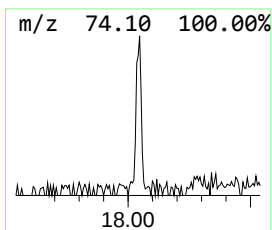
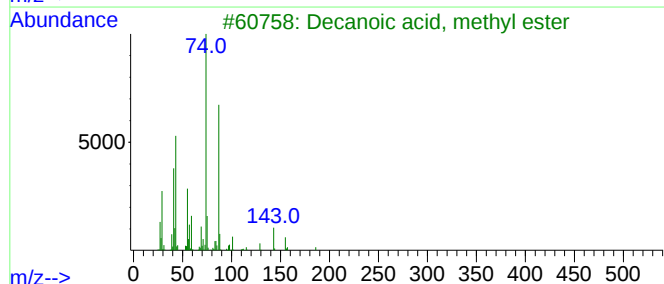
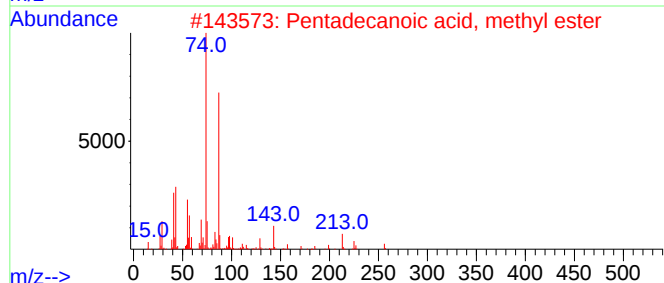
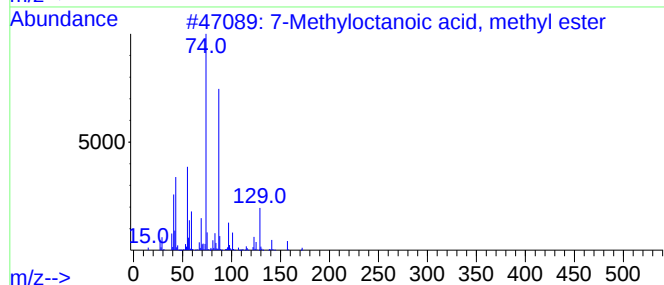
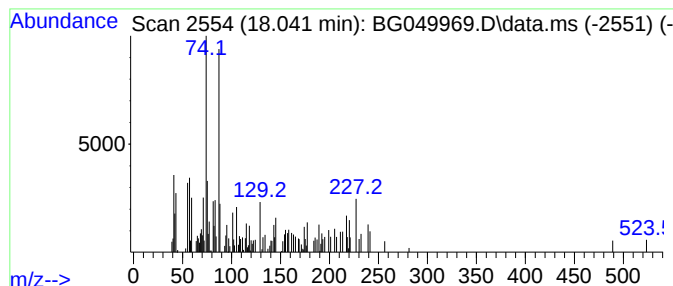
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 7-Methyloctanoic acid, meth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.041	2.32 ng	41350	Phenanthrene-d10	17.395

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Methyloctanoic acid, methyl ester	172	C10H20O2	005129-53-3	64
2			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	52
3			Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	50
4			Nonanoic acid, methyl ester	172	C10H20O2	001731-84-6	50
5			Pentanoic acid, 4-methyl-, methy...	130	C7H14O2	002412-80-8	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082521\
Data File : BG049969.D
Acq On : 25 Aug 2021 15:31
Operator : CG/JU
Sample : M3506-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown8.066	8.066	2.4	ng	19690	1	7.978	162068	20.0
Triethyl phosphate	9.447	2.5	ng	30305	2	10.798	242637	20.0
Octasiloxane, 1...	9.852	5.8	ng	69978	2	10.798	242637	20.0
Ethanol, 2-(2-b...	10.557	12.0	ng	145192	2	10.798	242637	20.0
Methyl 2-(2-(2-...	11.409	3.2	ng	39305	2	10.798	242637	20.0
Dipropylene gly...	11.702	4.5	ng	54428	2	10.798	242637	20.0
Hexaethylene gl...	11.761	4.1	ng	49268	2	10.798	242637	20.0
Hydrazine, 1,1-...	11.943	3.0	ng	36431	2	10.798	242637	20.0
unknown15.045	15.045	2.2	ng	35641	3	14.634	318087	20.0
7-Methyloctanoi...	18.041	2.3	ng	41350	4	17.395	356978	20.0