

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080824\
 Data File : BF138868.D
 Acq On : 08 Aug 2024 16:28
 Operator : RC/JU
 Sample : P3416-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MLS-15-106-121

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_F\Methods\8270-BF073024.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF138868.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.898	302	307	317	rBB	18796	29585	1.68%	0.337%
2	5.051	499	503	515	rVB	18603	27200	1.54%	0.310%
3	5.469	569	574	593	rBV	356596	534207	30.26%	6.093%
4	5.798	623	630	654	rBV	605927	942392	53.38%	10.748%
5	6.481	741	746	768	rVB	237898	335615	19.01%	3.828%
6	6.839	802	807	813	rBB	180948	223757	12.67%	2.552%
7	7.404	897	903	917	rBV	722928	987011	55.90%	11.257%
8	8.116	1019	1024	1043	rBB	230814	295806	16.75%	3.374%
9	8.333	1057	1061	1063	rBV	83042	103351	5.85%	1.179%
10	8.363	1063	1066	1077	rVB	91571	135981	7.70%	1.551%
11	9.192	1201	1207	1212	rBV	1391497	1765554	100.00%	20.137%
12	9.869	1317	1322	1330	rBV	264336	322908	18.29%	3.683%
13	10.663	1452	1457	1489	rBV	703541	925186	52.40%	10.552%
14	11.357	1570	1575	1584	rBV	255470	314823	17.83%	3.591%
15	12.939	1838	1844	1849	rBV	1173835	1486476	84.19%	16.954%
16	13.998	2019	2024	2032	rVB	132730	178185	10.09%	2.032%
17	15.457	2266	2272	2286	rBV2	94558	159634	9.04%	1.821%

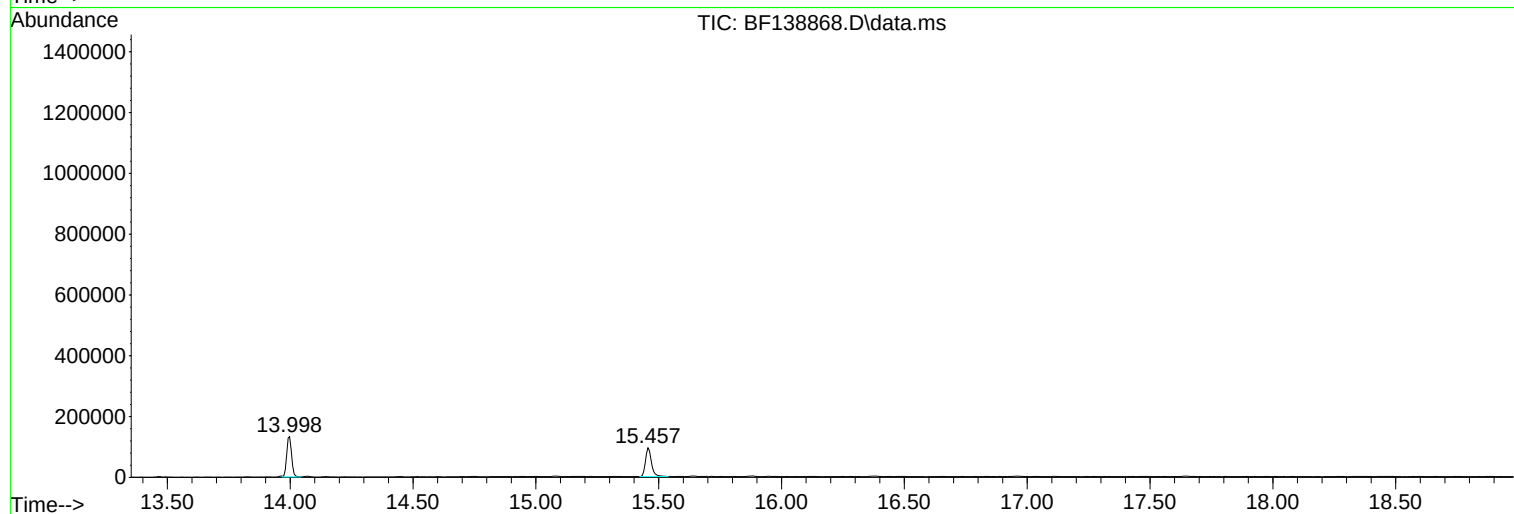
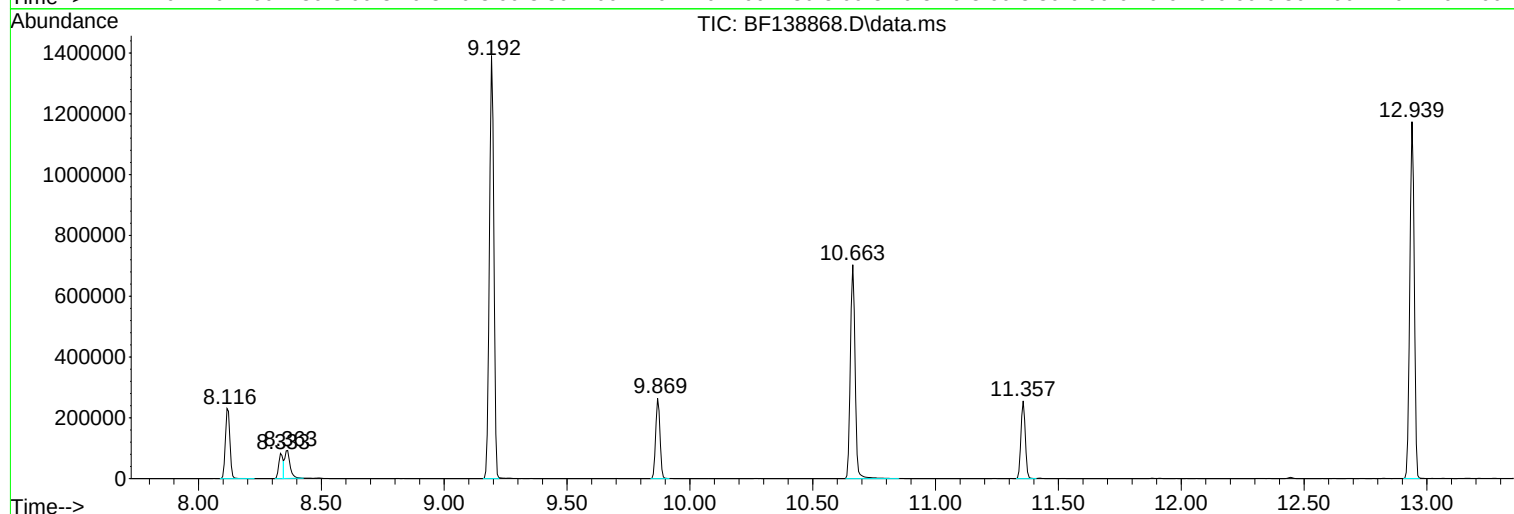
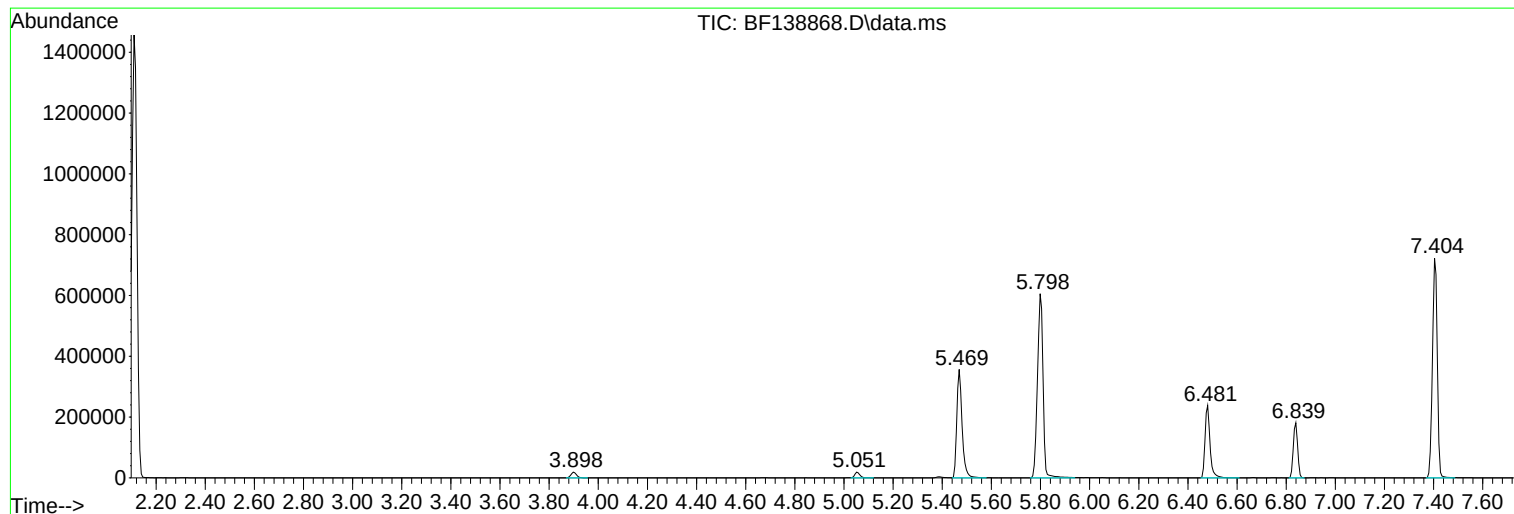
Sum of corrected areas: 8767671

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080824\
Data File : BF138868.D
Acq On : 08 Aug 2024 16:28
Operator : RC/JU
Sample : P3416-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MLS-15-106-121

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080824\
Data File : BF138868.D
Acq On : 08 Aug 2024 16:28
Operator : RC/JU
Sample : P3416-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MLS-15-106-121

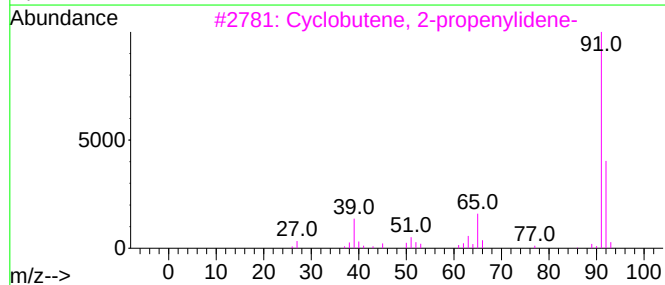
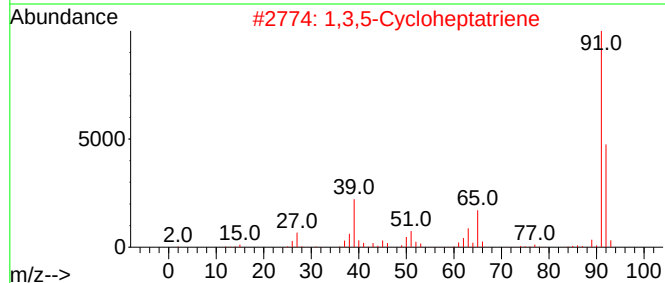
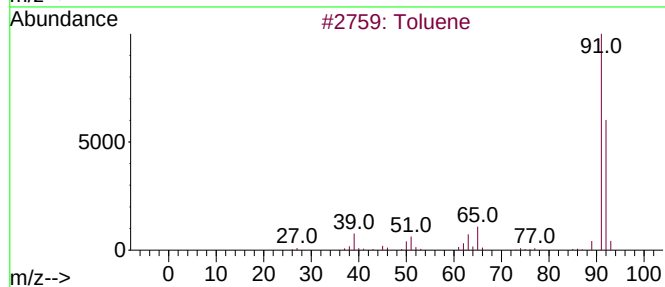
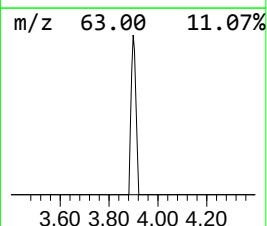
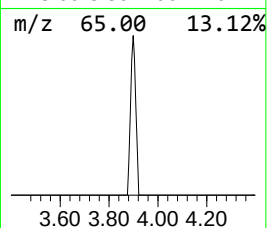
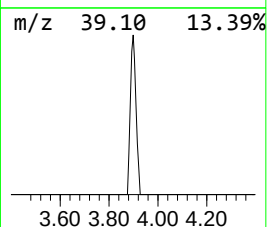
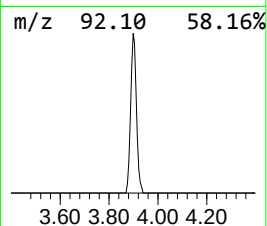
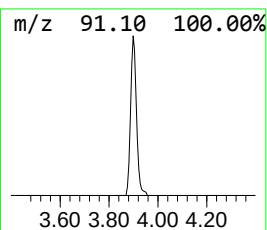
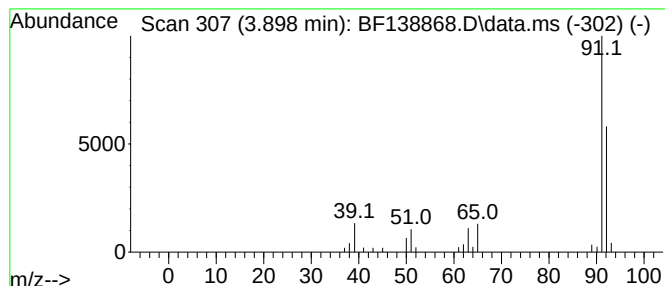
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Toluene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.898	2.64 ng	29585	1,4-Dichlorobenzene-d4	6.839

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Toluene	92	C7H8	000108-88-3	91
2			1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	86
3			Cyclobutene, 2-propenylidene-	92	C7H8	052097-85-5	80
4			1,5-Heptadien-3-yne	92	C7H8	003511-27-1	72
5			Molybdenum, di-.mu.-chlorobis[(1... 532	C20H26Cl2Mo2	035625-66-2	72	



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080824\
Data File : BF138868.D
Acq On : 08 Aug 2024 16:28
Operator : RC/JU
Sample : P3416-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MLS-15-106-121

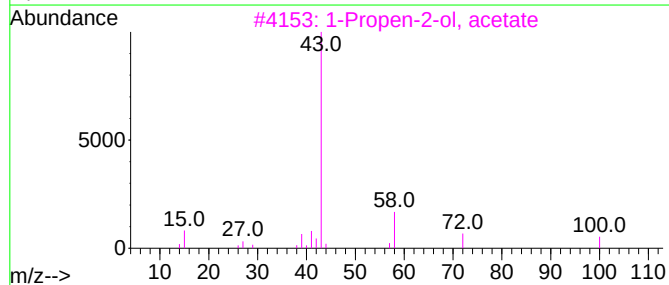
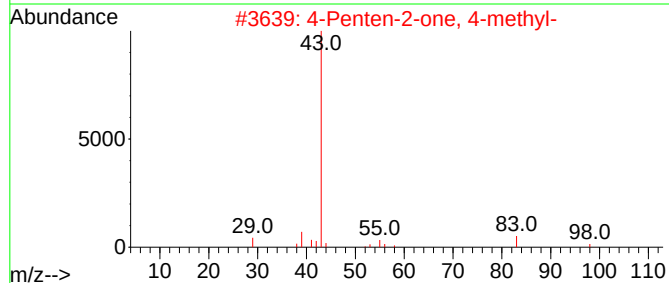
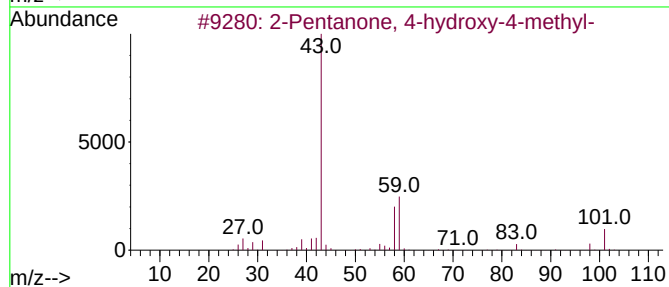
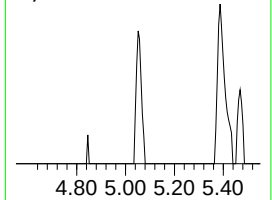
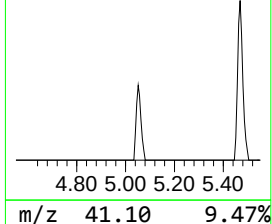
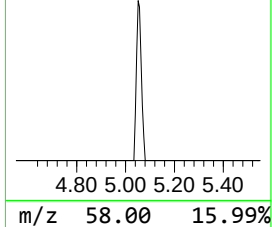
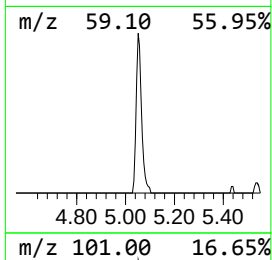
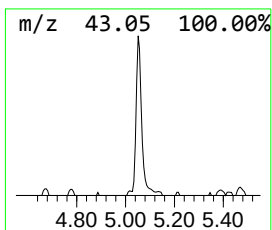
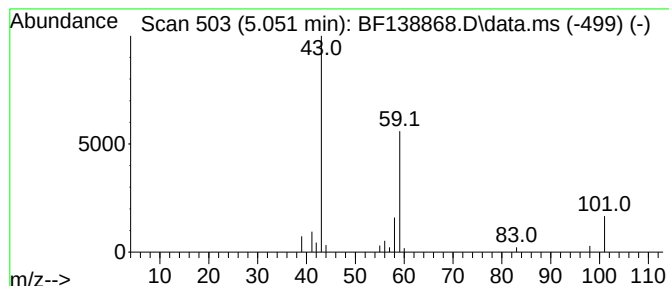
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.051	2.43 ng	27200	1,4-Dichlorobenzene-d4	6.839

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
4		3-Hexanol	102	C6H14O	000623-37-0	9
5		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080824\
Data File : BF138868.D
Acq On : 08 Aug 2024 16:28
Operator : RC/JU
Sample : P3416-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MLS-15-106-121

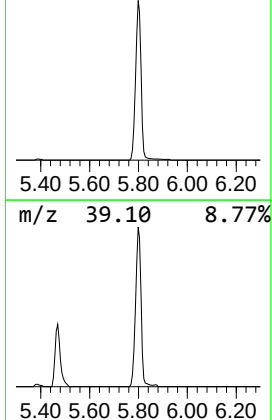
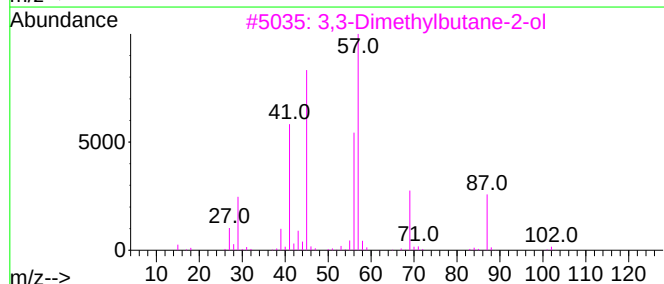
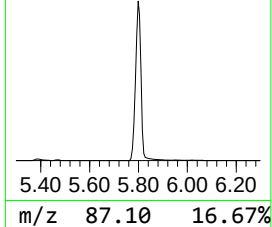
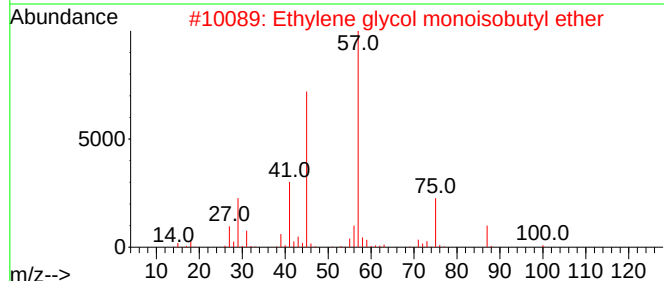
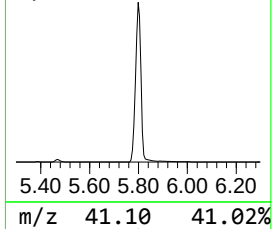
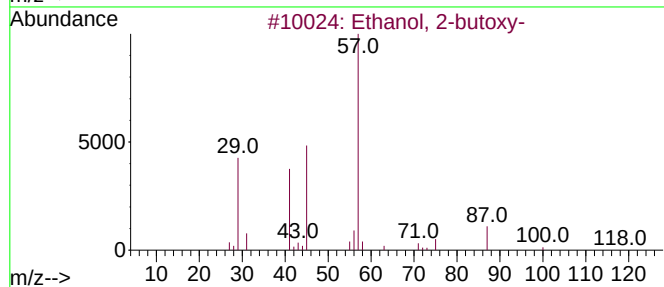
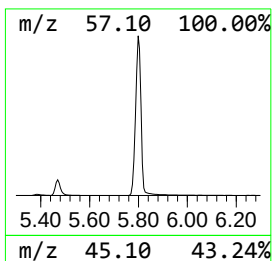
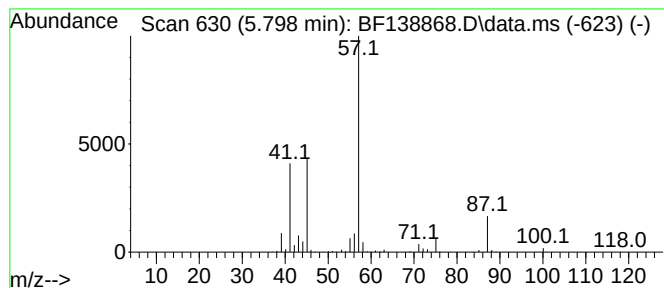
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Ethanol, 2-butoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.798	84.23 ng	942392	1,4-Dichlorobenzene-d4	6.839

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	83
2			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	72
3			3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	42
4			Propanoic acid, 2,2-dimethyl-	102	C5H10O2	000075-98-9	25
5			1,3-Propanediol, 2-methyl-, dipr...	202	C10H18O4	054932-83-1	17



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080824\
Data File : BF138868.D
Acq On : 08 Aug 2024 16:28
Operator : RC/JU
Sample : P3416-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MLS-15-106-121

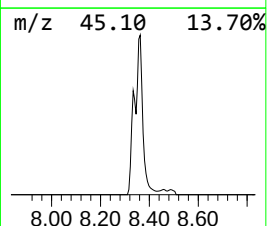
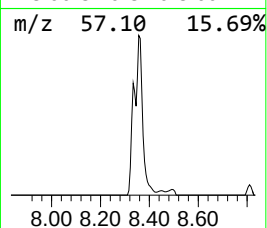
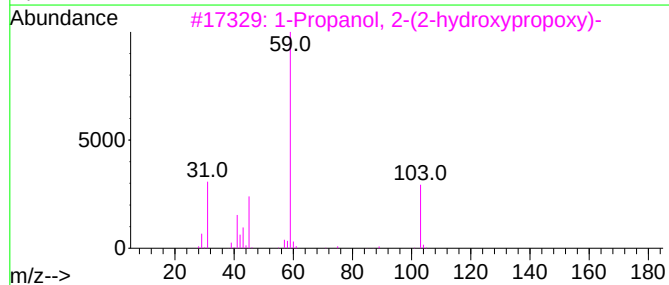
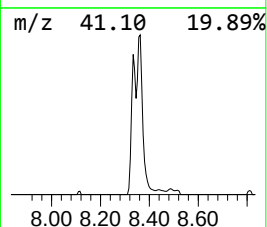
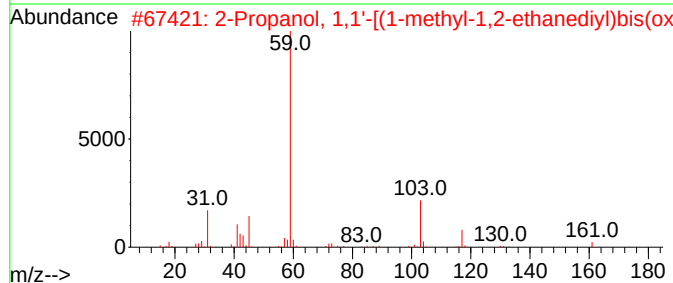
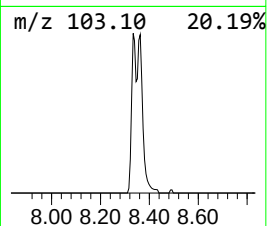
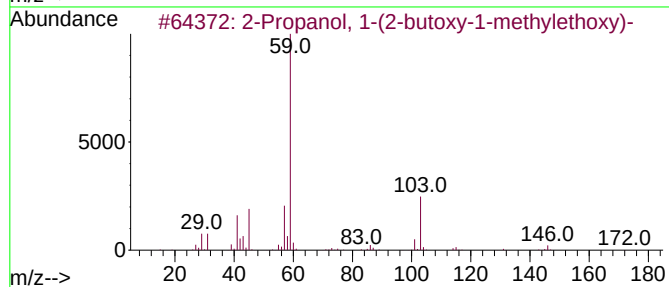
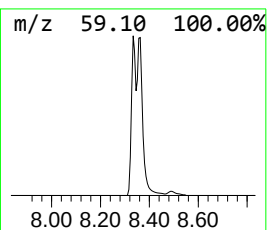
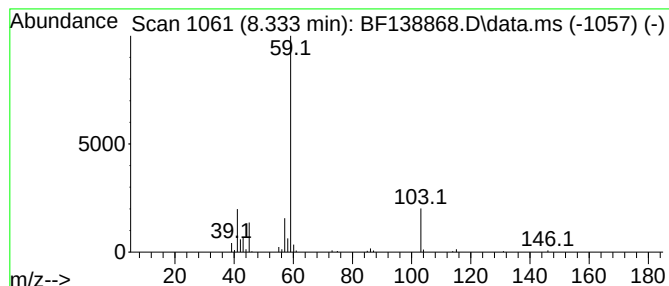
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 2-Propanol, 1-(2-butoxy-1-m... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.333	6.99 ng	103351	Naphthalene-d8	8.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	83		
2	2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	78		
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78		
4	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	74		
5	Tetraglyme	222	C10H22O5	000143-24-8	72		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080824\
Data File : BF138868.D
Acq On : 08 Aug 2024 16:28
Operator : RC/JU
Sample : P3416-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MLS-15-106-121

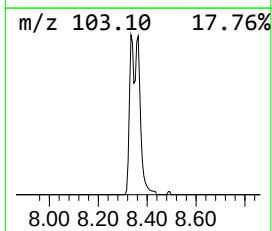
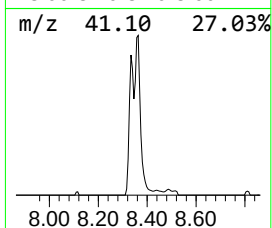
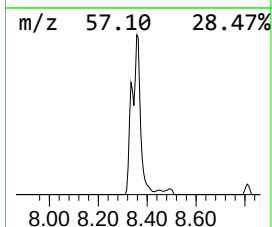
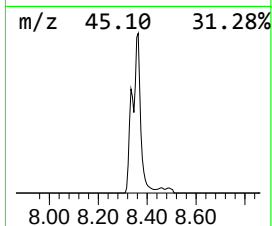
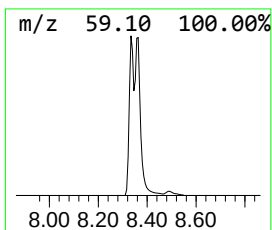
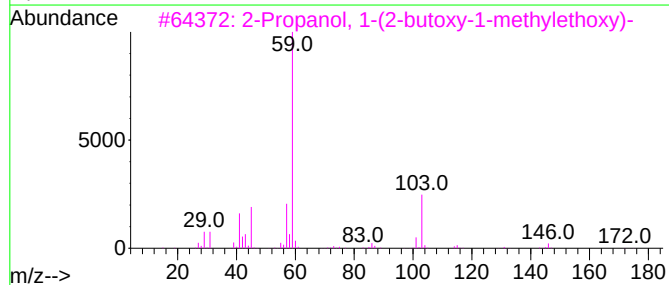
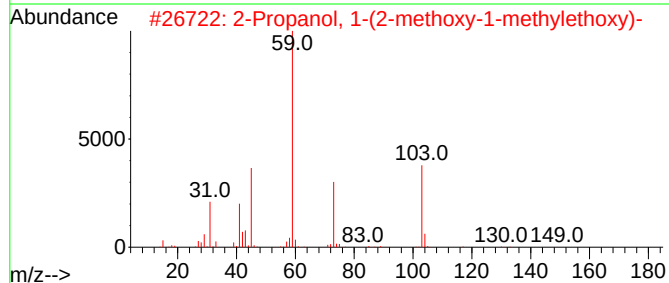
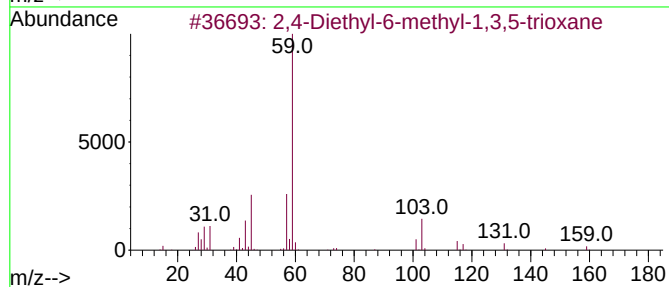
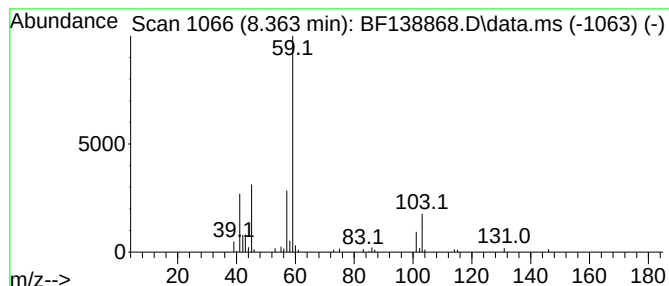
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 2,4-Diethyl-6-methyl-1,3,5-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.363	9.19 ng	135981	Naphthalene-d8	8.116

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Diethyl-6-methyl-1,3,5-trioxane	160	C8H16O3	117888-04-7	72	
2	2-Propanol, 1-(2-methoxy-1-methyl-2-propoxy)-	148	C7H16O3	020324-32-7	64	
3	2-Propanol, 1-(2-butoxy-1-methyl-2-propoxy)-	190	C10H22O3	029911-28-2	64	
4	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	59	
5	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	59	



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080824\
Data File : BF138868.D
Acq On : 08 Aug 2024 16:28
Operator : RC/JU
Sample : P3416-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MLS-15-106-121

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Toluene	3.898	2.6	ng	29585	1	6.839	223757	20.0
2-Pentanone, 4-...	5.051	2.4	ng	27200	1	6.839	223757	20.0
Ethanol, 2-butoxy-	5.798	84.2	ng	942392	1	6.839	223757	20.0
2-Propanol, 1-(...	8.333	7.0	ng	103351	2	8.116	295806	20.0
2,4-Diethyl-6-m...	8.363	9.2	ng	135981	2	8.116	295806	20.0