

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042722\
 Data File : BF127959.D
 Acq On : 27 Apr 2022 21:50
 Operator : CG\JU
 Sample : N2512-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-19

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-BF042222.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF127959.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.251	328	334	338	rBV	53716	65699	1.70%	0.322%
2	4.457	365	369	377	rBV	258733	279324	7.24%	1.369%
3	5.157	484	488	495	rBV	123948	136750	3.54%	0.670%
4	5.545	550	554	568	rVB	1600351	1389250	35.99%	6.809%
5	6.545	720	724	730	rVB	929061	874074	22.65%	4.284%
6	6.728	752	755	757	rBV	158528	161449	4.18%	0.791%
7	6.757	757	760	769	rVV2	184064	290463	7.53%	1.424%
8	6.851	773	776	786	rVV	203035	302286	7.83%	1.482%
9	6.928	786	789	793	rVB	777231	653716	16.94%	3.204%
10	7.063	801	812	813	rBV3	27553	72125	1.87%	0.354%
11	7.263	843	846	853	rVB2	39733	55537	1.44%	0.272%
12	7.334	853	858	866	rVB	72913	80982	2.10%	0.397%
13	7.492	881	885	888	rVB	2110277	2175991	56.38%	10.666%
14	7.575	894	899	902	rBV	68393	81440	2.11%	0.399%
15	7.633	906	909	917	rVB4	19106	41454	1.07%	0.203%
16	7.928	955	959	965	rVB3	48962	56338	1.46%	0.276%
17	8.175	992	1001	1003	rBV5	30357	74488	1.93%	0.365%
18	8.210	1003	1007	1011	rVV	895349	823803	21.34%	4.038%
19	8.439	1039	1046	1047	rBV2	25424	38737	1.00%	0.190%
20	8.545	1060	1064	1069	rBV2	97820	115675	3.00%	0.567%
21	8.616	1073	1076	1087	rVB	47678	71174	1.84%	0.349%
22	9.204	1172	1176	1178	rBV3	49658	62619	1.62%	0.307%
23	9.292	1185	1191	1194	rBV	3729733	3859719	100.00%	18.918%
24	9.386	1204	1207	1223	rVB	558417	768321	19.91%	3.766%
25	9.969	1299	1306	1310	rBV	1034668	920004	23.84%	4.509%
26	10.380	1372	1376	1379	rVB	190660	157823	4.09%	0.774%
27	10.674	1422	1426	1428	rBV	123623	105489	2.73%	0.517%
28	10.763	1436	1441	1444	rBV	2188686	1938875	50.23%	9.503%
29	11.463	1556	1560	1563	rBV	983472	817383	21.18%	4.006%
30	11.839	1621	1624	1627	rVB	69336	50120	1.30%	0.246%
31	13.051	1825	1830	1833	rBV	3021375	2664034	69.02%	13.058%
32	14.104	2006	2009	2013	rBV	581139	539992	13.99%	2.647%
33	15.615	2260	2266	2274	rBV	512854	676837	17.54%	3.318%

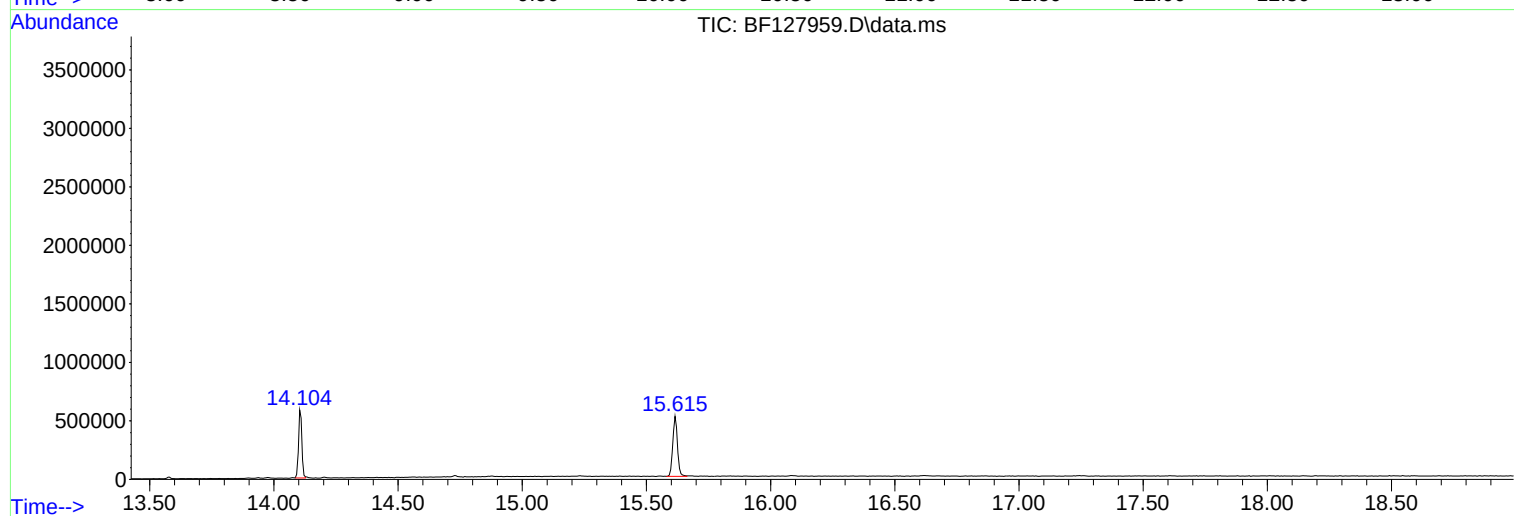
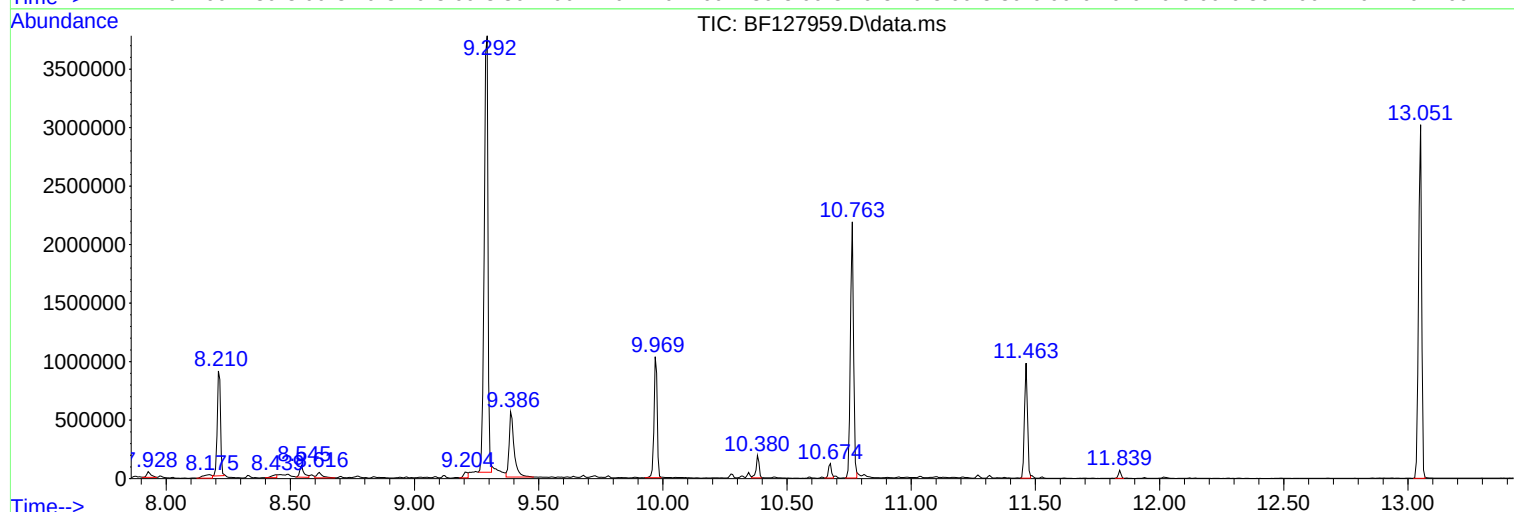
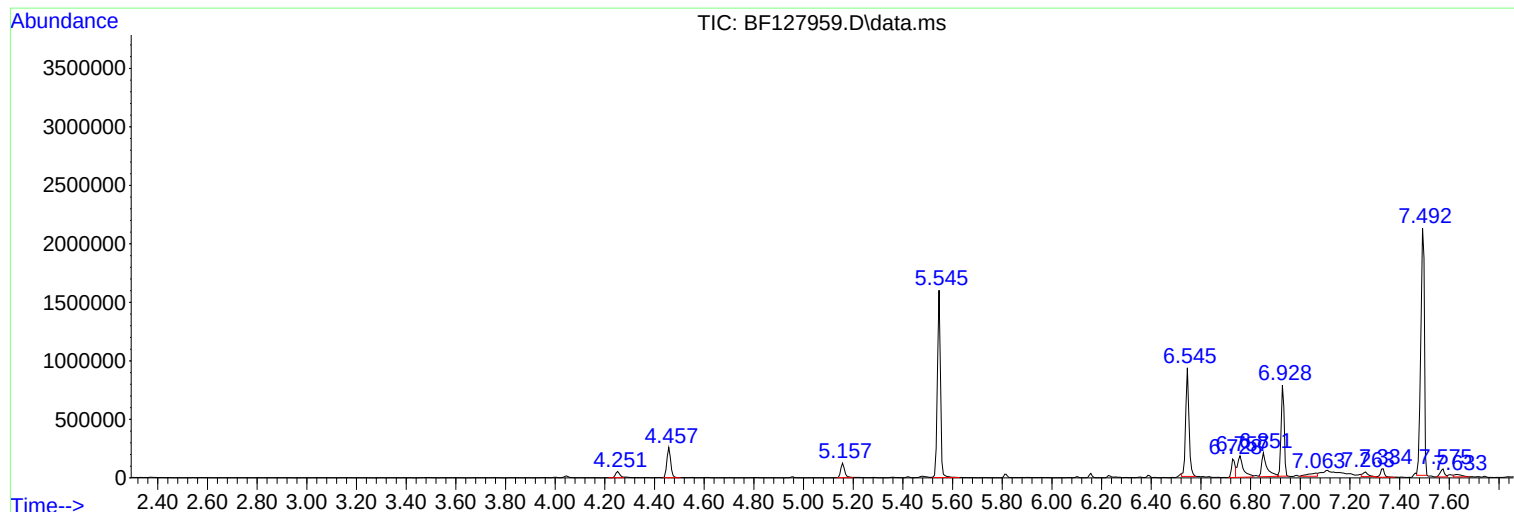
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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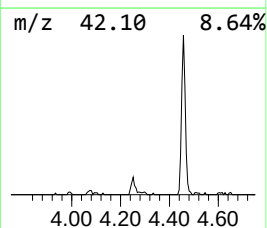
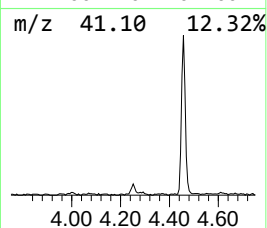
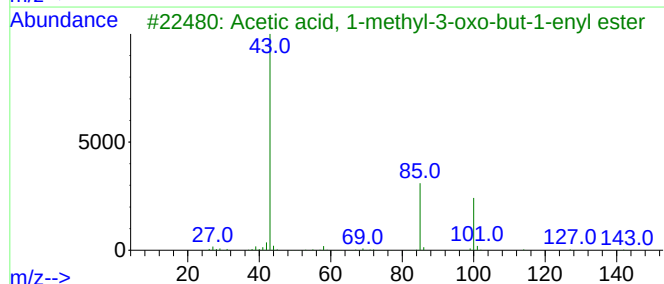
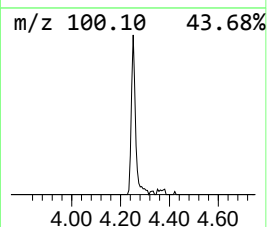
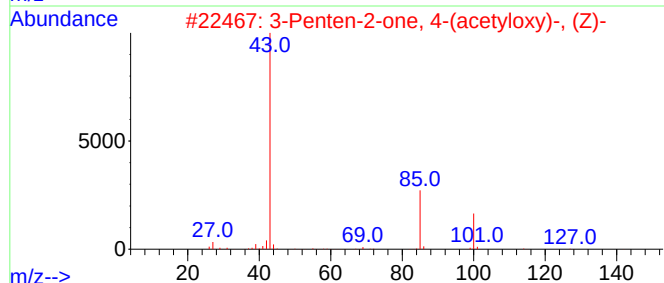
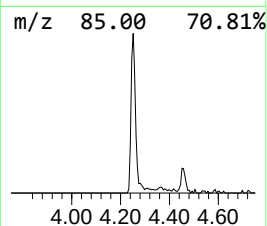
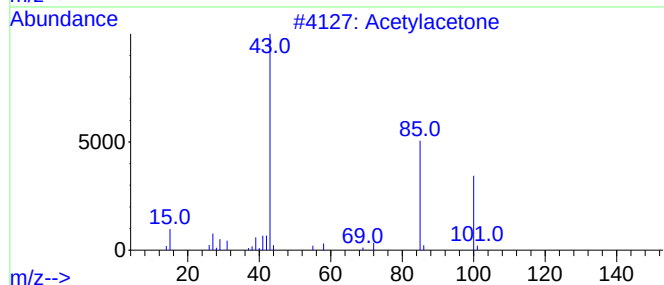
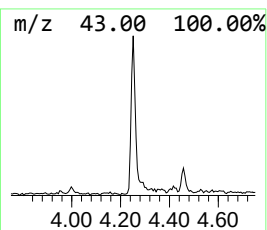
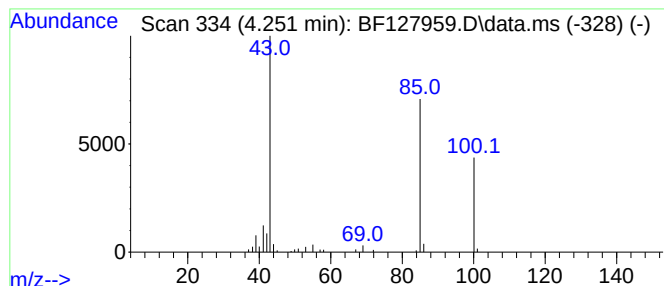
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Acetylacetone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.251	2.01 ng	65699	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetylacetone	100	C5H8O2	000123-54-6	90
2			3-Penten-2-one, 4-(acetyloxy)-, ...	142	C7H10O3	038365-58-1	43
3			Acetic acid, 1-methyl-3-oxo-but-...	142	C7H10O3	1000190-85-2	43
4			Hexane, 3-bromo-	164	C6H13Br	003377-87-5	43
5			Hexane, 2-bromo-	164	C6H13Br	003377-86-4	43



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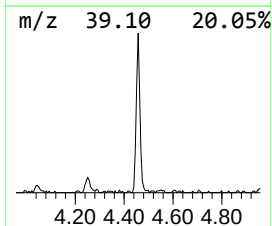
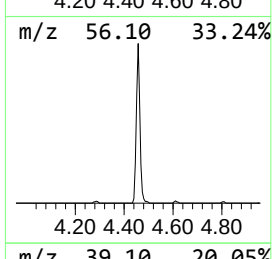
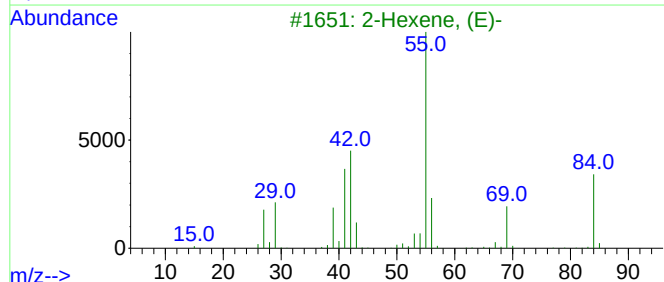
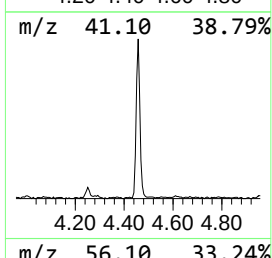
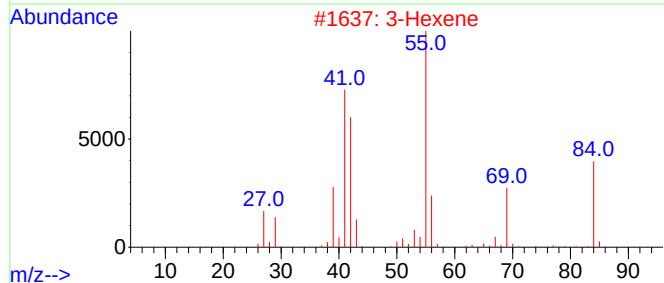
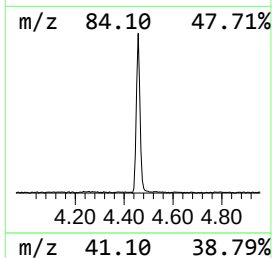
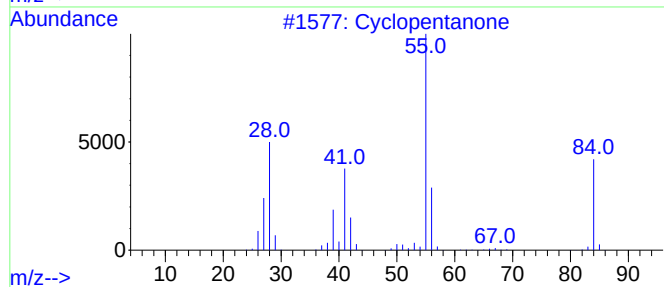
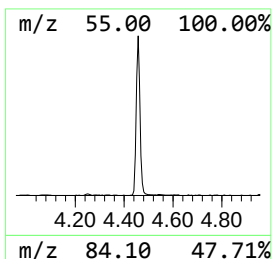
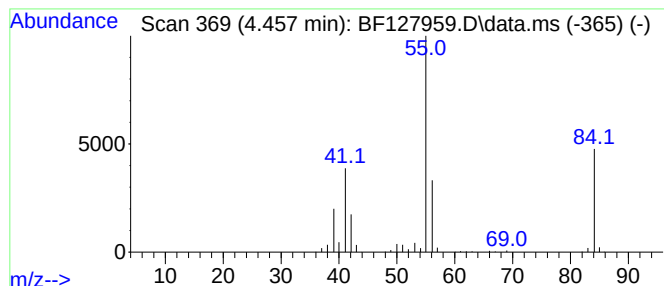
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclopentanone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.457	8.55 ng	279324	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentanone	84	C5H8O	000120-92-3	90
2			3-Hexene	84	C6H12	000592-47-2	72
3			2-Hexene, (E)-	84	C6H12	004050-45-7	64
4			2-Hexene	84	C6H12	000592-43-8	64
5			3-Hexene, (Z)-	84	C6H12	007642-09-3	50



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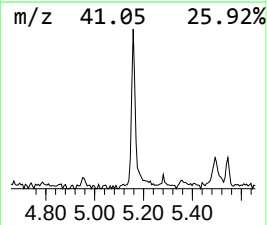
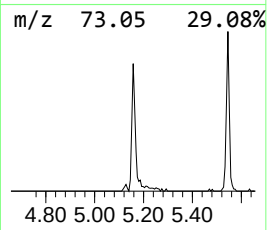
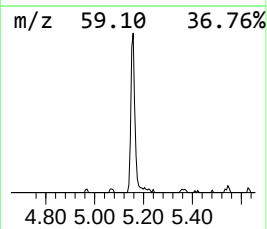
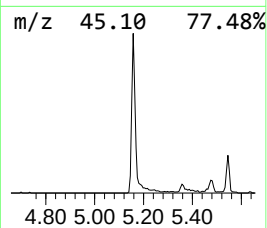
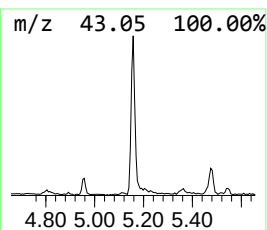
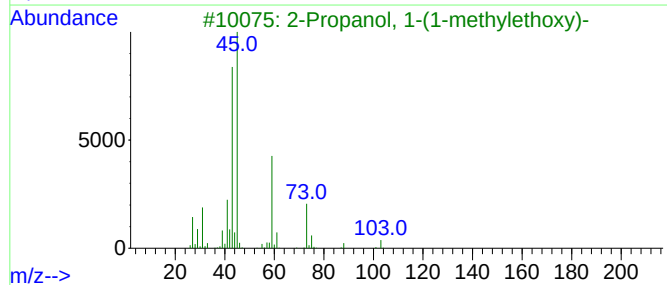
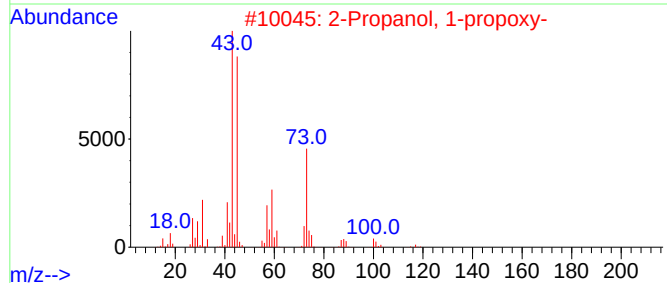
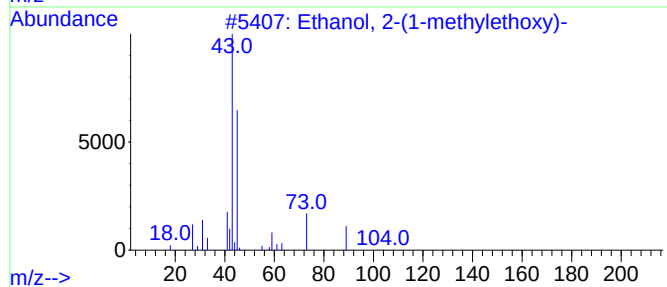
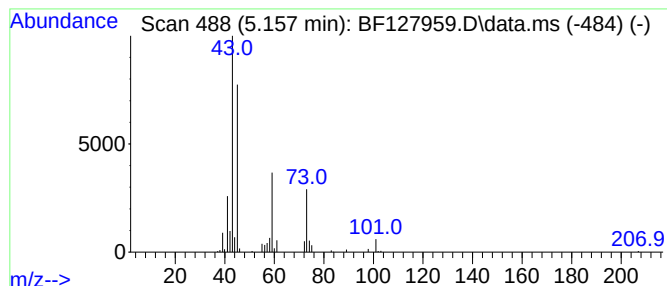
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Ethanol, 2-(1-methylethoxy)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.157	4.18 ng	136750	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(1-methylethoxy)-	104	C5H12O2	000109-59-1	72
2			2-Propanol, 1-propoxy-	118	C6H14O2	001569-01-3	64
3			2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	59
4			Butane, 1-(1-methylethoxy)-	116	C7H16O	001860-27-1	50
5			Sulfurous acid, bis(1-methylethy...	166	C6H14O3S	004773-13-1	45



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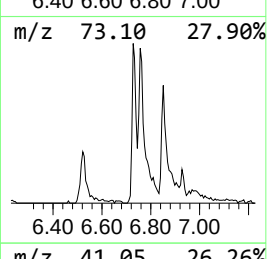
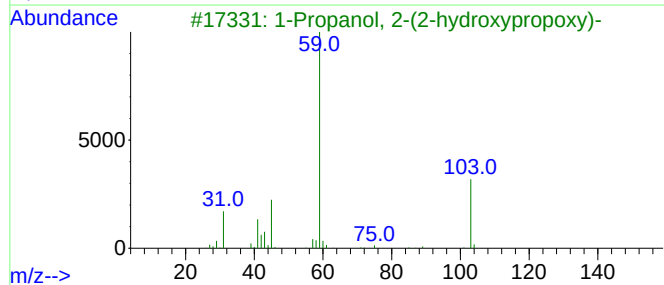
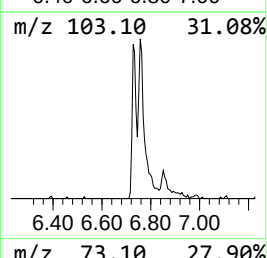
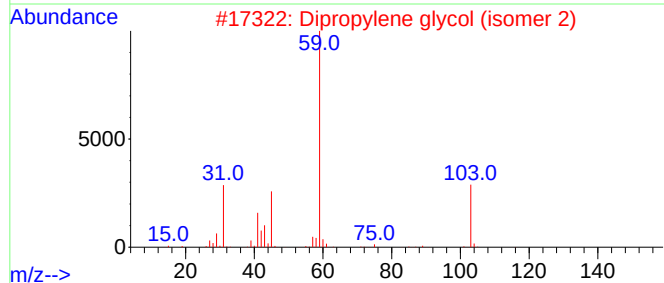
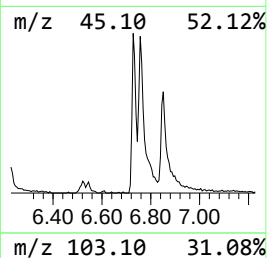
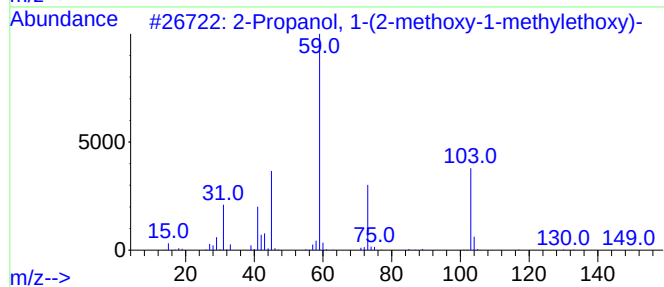
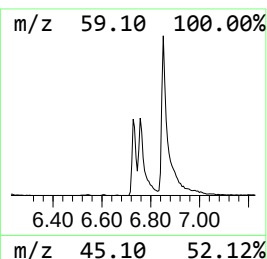
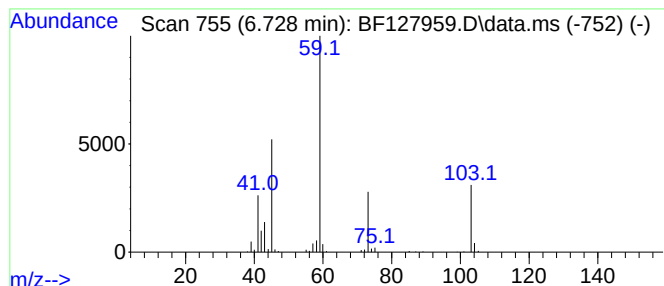
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.728	4.94 ng	161449	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	86
2			Dipropylene glycol (isomer 2)	134	C6H14O3	1000506-25-4	59
3			1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	59
4			Ethane, 1-ethoxy-1-methoxy-	104	C5H12O2	010471-14-4	59
5			Dipropylene glycol (isomer 3)	134	C6H14O3	1000506-25-5	53



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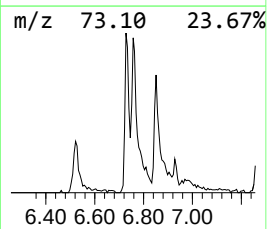
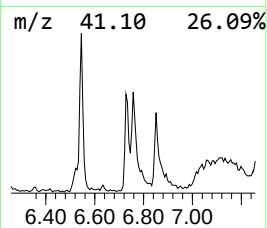
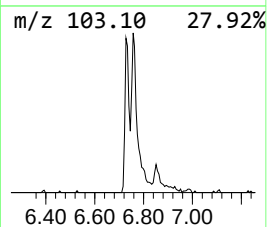
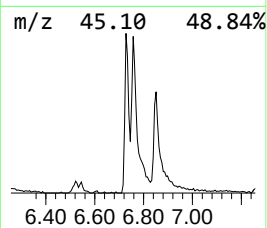
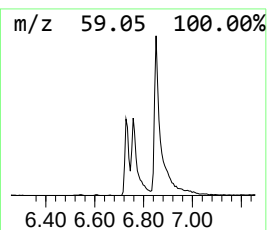
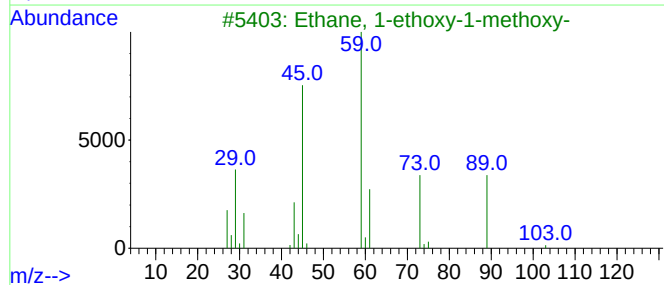
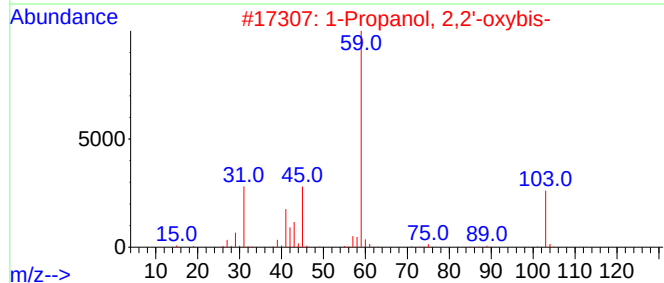
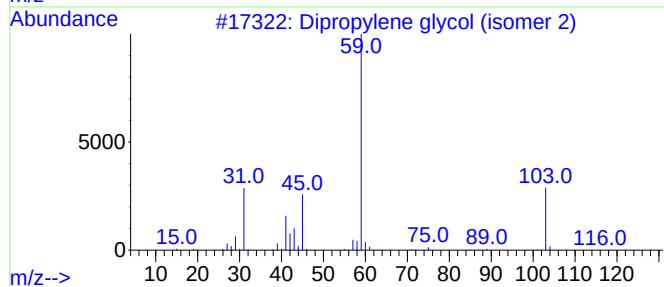
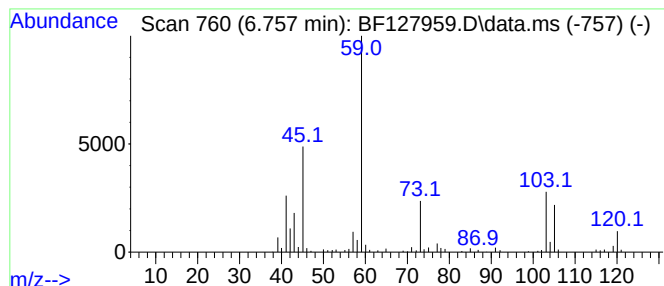
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Peak Number 5 Dipropylene glycol (isomer 2) Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.757	8.89 ng	290463	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dipropylene glycol (isomer 2)	134	C6H14O3	1000506-25-4	50
2			1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	50
3			Ethane, 1-ethoxy-1-methoxy-	104	C5H12O2	010471-14-4	50
4			Dipropylene glycol (isomer 3)	134	C6H14O3	1000506-25-5	50
5			2-Propanol, 1-(2-ethoxypropoxy)-	162	C8H18O3	010143-32-5	45



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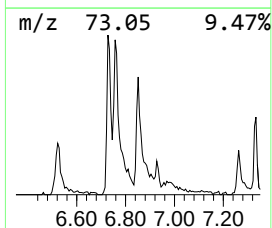
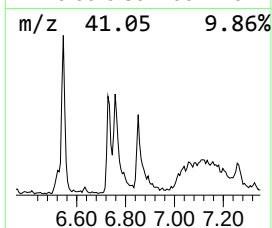
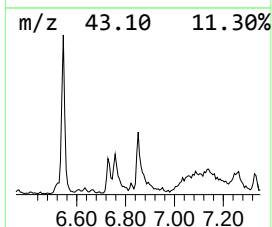
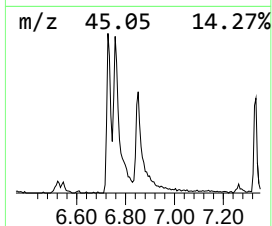
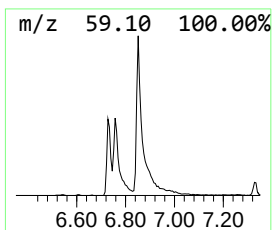
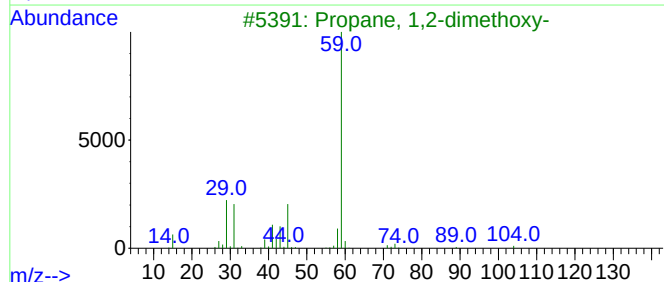
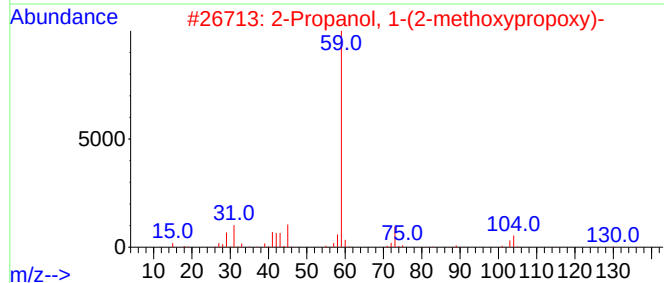
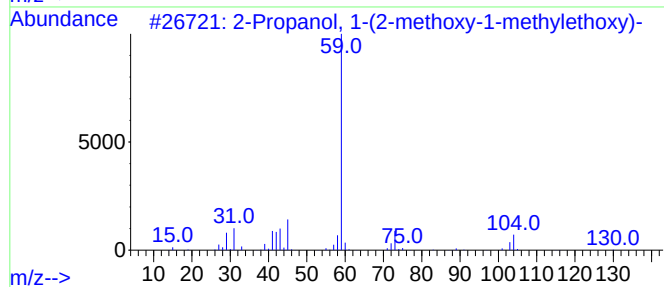
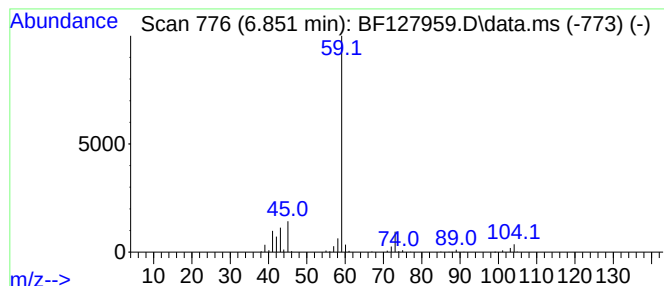
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2-Propanol, 1-(2-methoxypro... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.851	9.25 ng	302286	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	83
2			2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	83
3			Propane, 1,2-dimethoxy-	104	C5H12O2	007778-85-0	72
4			1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64
5			Tetraglyme	222	C10H22O5	000143-24-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042722\
Data File : BF127959.D
Acq On : 27 Apr 2022 21:50
Operator : CG\JU
Sample : N2512-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-19

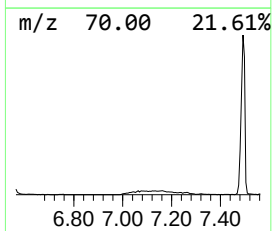
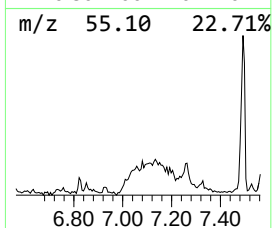
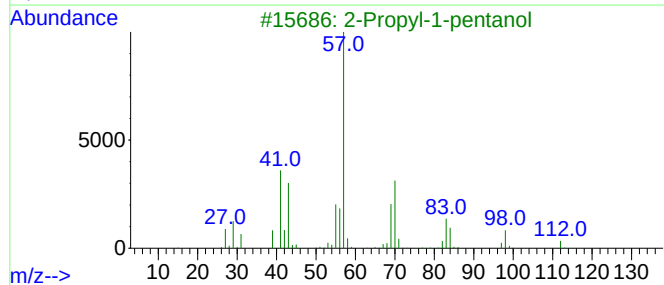
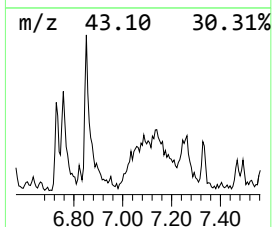
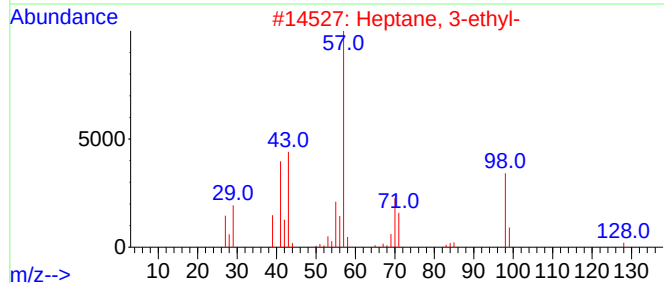
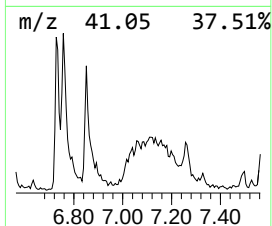
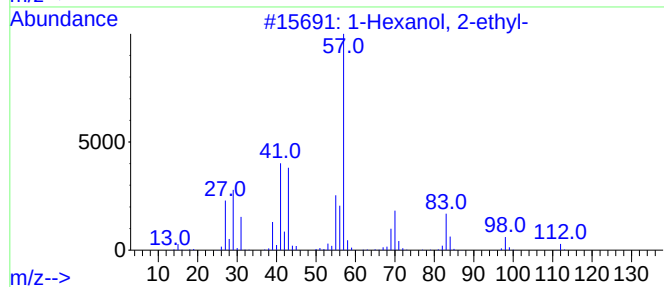
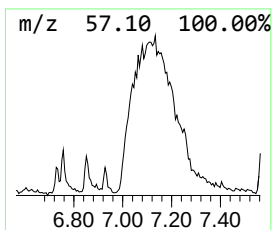
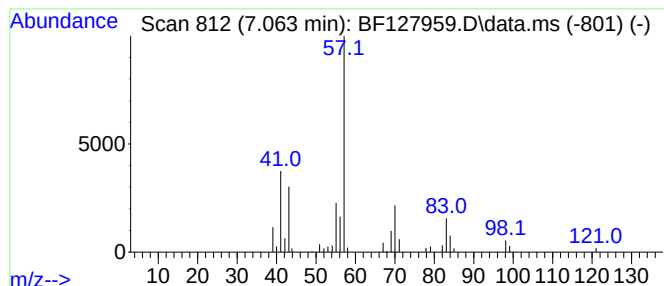
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1-Hexanol, 2-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.063	2.21 ng	72125	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2			Heptane, 3-ethyl-	128	C9H20	015869-80-4	50
3			2-Propyl-1-pentanol	130	C8H18O	058175-57-8	50
4			1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	50
5			2-Ethyl-1-hexanol, trifluoroacetate	226	C10H17F3O2	053800-08-1	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042722\
Data File : BF127959.D
Acq On : 27 Apr 2022 21:50
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Sample : N2512-01
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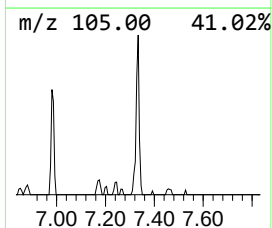
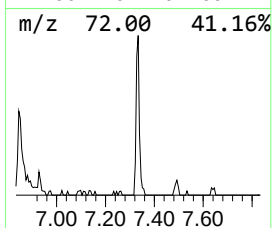
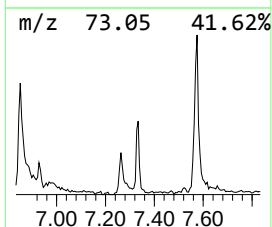
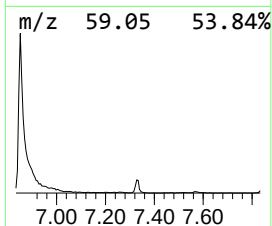
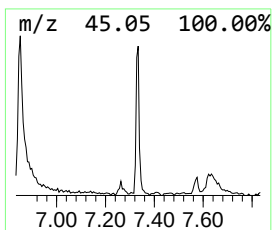
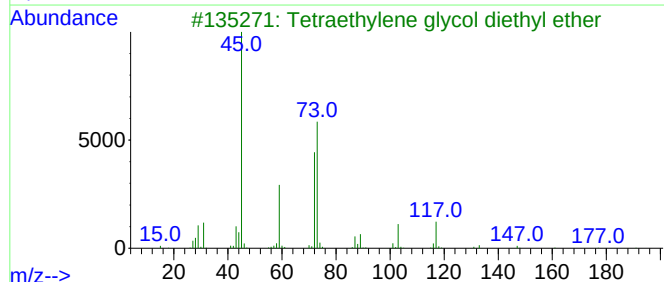
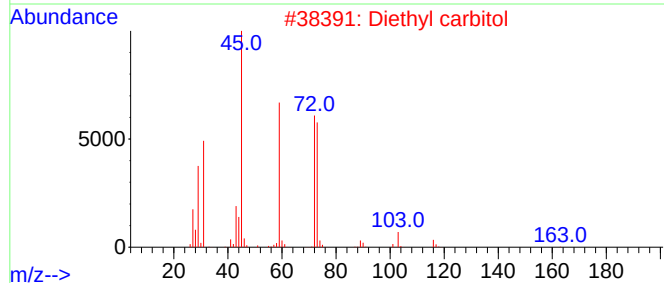
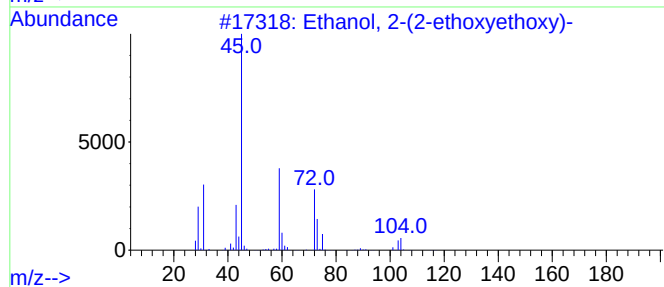
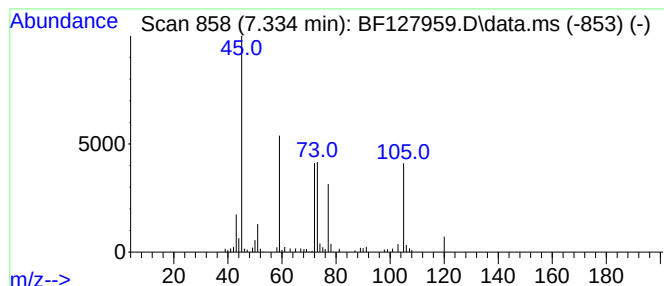
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.334	2.48 ng	80982	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	59
2			Diethyl carbitol	162	C8H18O3	000112-36-7	45
3			Tetraethylene glycol diethyl ether	250	C12H26O5	004353-28-0	33
4			Ethanol, 2-[2-(2-methoxyethoxy)e...	164	C7H16O4	000112-35-6	32
5			Propane, 1,3-dimethoxy-	104	C5H12O2	017081-21-9	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042722\
Data File : BF127959.D
Acq On : 27 Apr 2022 21:50
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Sample : N2512-01
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ClientSampleId :
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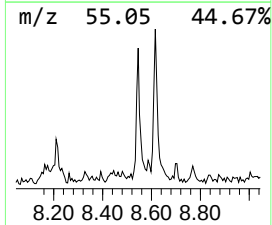
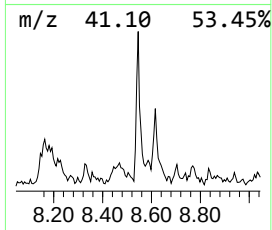
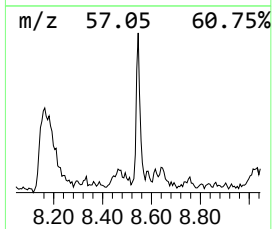
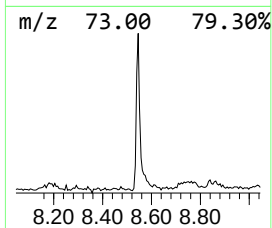
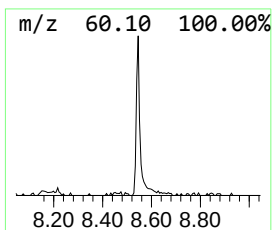
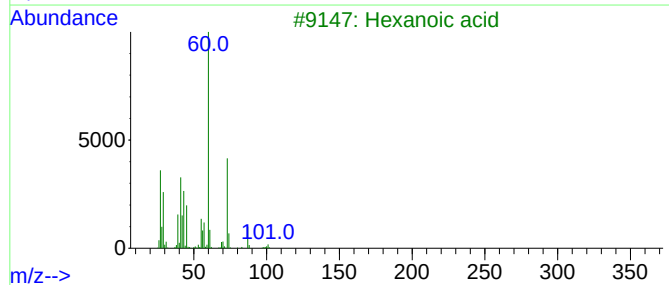
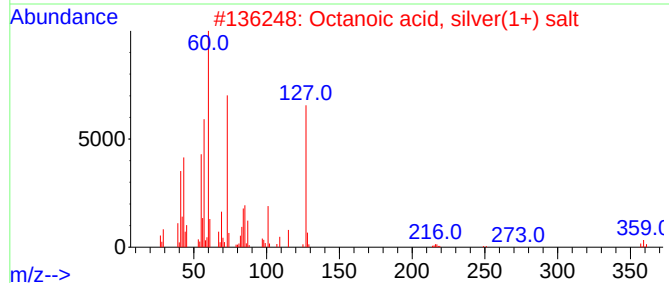
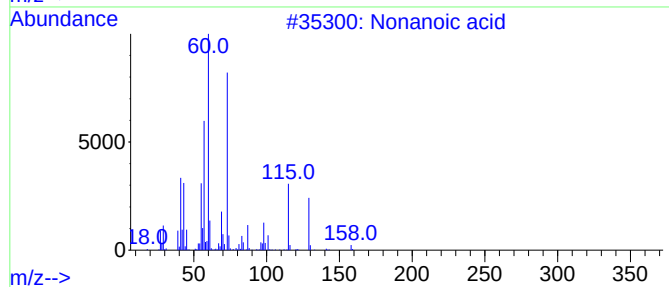
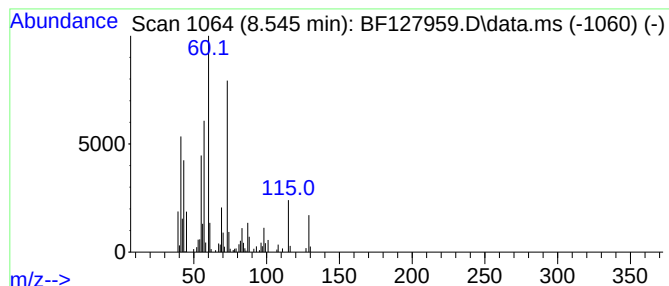
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Nonanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.545	2.81 ng	115675	Naphthalene-d8	8.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonanoic acid	158	C9H18O2	000112-05-0	90
2			Octanoic acid, silver(1+) salt	250	C8H15AgO2	024927-67-1	50
3			Hexanoic acid	116	C6H12O2	000142-62-1	50
4			Octanoic acid	144	C8H16O2	000124-07-2	45
5			Propanedioic acid, propyl-	146	C6H10O4	000616-62-6	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042722\
Data File : BF127959.D
Acq On : 27 Apr 2022 21:50
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Sample : N2512-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

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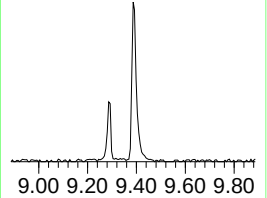
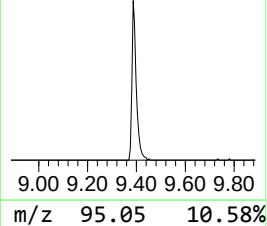
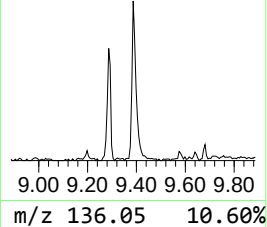
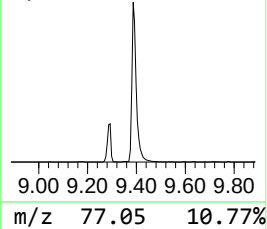
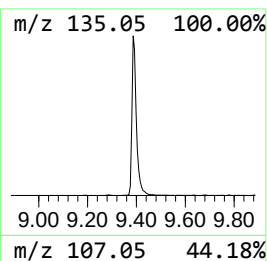
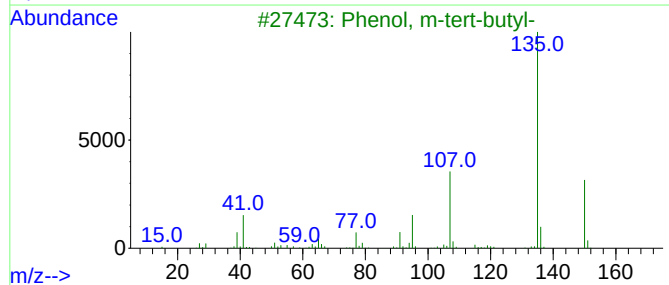
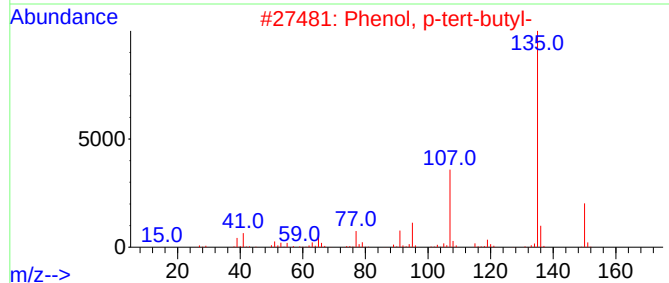
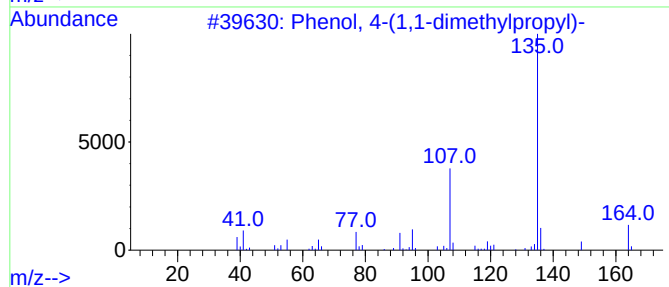
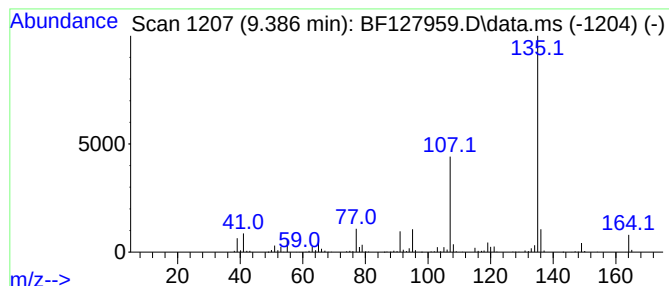
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenol, 4-(1,1-dimethylprop... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.386	16.70 ng	768321	Acenaphthene-d10	9.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	93
2			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	83
3			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	64
4			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	64
5			Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042722\
Data File : BF127959.D
Acq On : 27 Apr 2022 21:50
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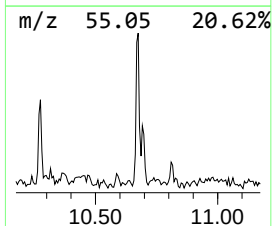
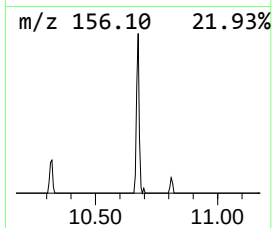
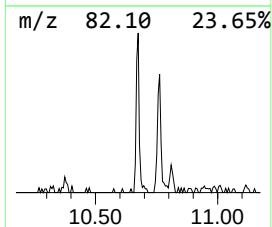
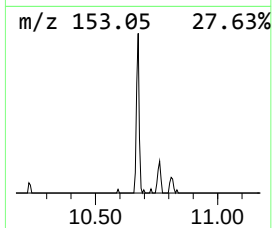
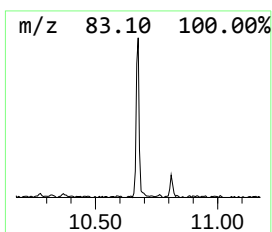
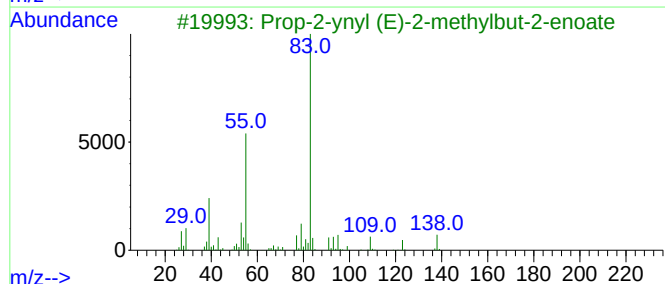
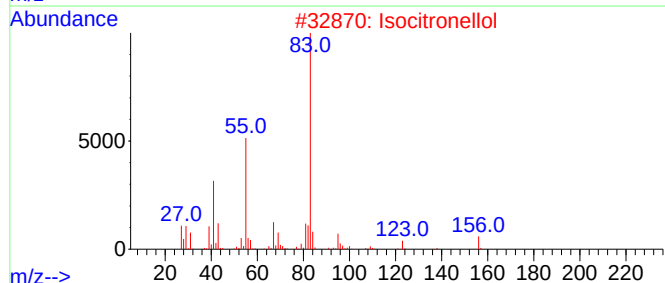
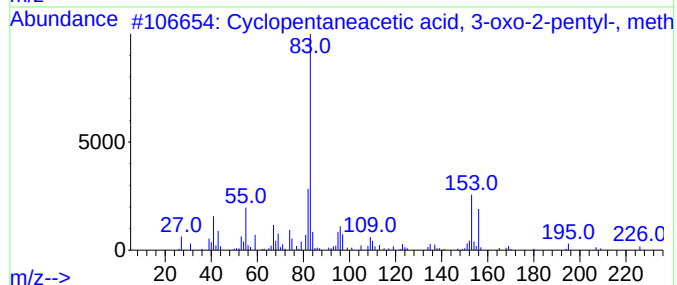
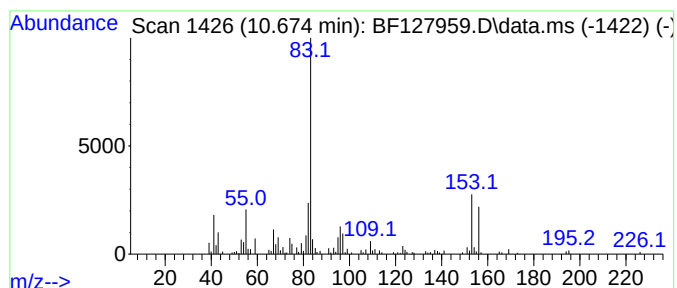
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Cyclopentaneacetic acid, 3-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.675	2.29 ng	105489	Acenaphthene-d10	9.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentaneacetic acid, 3-oxo-2...	226	C13H22O3	024851-98-7	90
2			Isocitronellol	156	C10H20O	018479-52-2	50
3			Prop-2-ynyl (E)-2-methylbut-2-en...	138	C8H10O2	1000373-72-5	43
4			Undecenyl tiglate, 2E-	252	C16H28O2	1000383-56-5	43
5			1,5-Dicyano-2,4-dimethyl-2,4-dia...	152	C7H12N4	1010289-23-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042722\
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 Acq On : 27 Apr 2022 21:50
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 Sample : N2512-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042222.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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Cyclopentanone	4.457	8.6	ng	279324	1	6.928	653716	20.0
Ethanol, 2-(1-m...	5.157	4.2	ng	136750	1	6.928	653716	20.0
2-Propanol, 1-(...	6.728	4.9	ng	161449	1	6.928	653716	20.0
Dipropylene gly...	6.757	8.9	ng	290463	1	6.928	653716	20.0
2-Propanol, 1-(...	6.851	9.3	ng	302286	1	6.928	653716	20.0
1-Hexanol, 2-et...	7.063	2.2	ng	72125	1	6.928	653716	20.0
Ethanol, 2-(2-e...	7.334	2.5	ng	80982	1	6.928	653716	20.0
Nonanoic acid	8.545	2.8	ng	115675	2	8.210	823803	20.0
Phenol, 4-(1,1-...	9.386	16.7	ng	768321	3	9.969	920004	20.0
Cyclopentaneace...	10.675	2.3	ng	105489	3	9.969	920004	20.0