

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072621\
 Data File : BF124777.D
 Acq On : 26 Jul 2021 22:05
 Operator : JU/CG
 Sample : M3151-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-1

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-BF072221.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF124777.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.163	10	13	17	rBB	9435	10190	7.83%	1.104%
2	5.099	507	512	515	rBB	8537	10590	8.14%	1.147%
3	5.493	575	579	582	rBB	135291	116194	89.31%	12.588%
4	6.498	745	750	753	rBV	126614	116598	89.63%	12.632%
5	6.875	811	814	817	rBB	38368	27893	21.44%	3.022%
6	7.434	905	909	912	rBB	80934	74374	57.17%	8.057%
7	8.051	1012	1014	1018	rBB	7430	5910	4.54%	0.640%
8	8.157	1029	1032	1035	rBB	44227	34882	26.81%	3.779%
9	9.234	1211	1215	1218	rBV	153826	130095	100.00%	14.094%
10	9.910	1327	1330	1333	rBB	47788	34521	26.54%	3.740%
11	10.698	1461	1464	1467	rBB	81470	61014	46.90%	6.610%
12	11.398	1580	1583	1586	rBB	39277	29516	22.69%	3.198%
13	11.916	1669	1671	1673	rBB	5268	3586	2.76%	0.388%
14	12.704	1803	1805	1807	rBB	2833	1790	1.38%	0.194%
15	12.986	1849	1853	1856	rBB	133978	113018	86.87%	12.244%
16	13.551	1946	1949	1951	rBB	4935	3921	3.01%	0.425%
17	13.880	2002	2005	2014	rBB3	10788	13660	10.50%	1.480%
18	14.039	2029	2032	2035	rBB	33228	25847	19.87%	2.800%
19	14.198	2056	2059	2062	rBV	11835	10048	7.72%	1.089%
20	14.498	2106	2110	2114	rBV	11786	12136	9.33%	1.315%
21	14.698	2138	2144	2146	rVB2	1637	2246	1.73%	0.243%
22	14.792	2156	2160	2162	rBV	17600	18149	13.95%	1.966%
23	14.810	2162	2163	2169	rVB	12919	9742	7.49%	1.055%
24	14.886	2174	2176	2178	rVB2	2647	2022	1.55%	0.219%
25	15.068	2201	2207	2209	rVB2	1152	1991	1.53%	0.216%
26	15.151	2217	2221	2225	rVB	10039	10774	8.28%	1.167%
27	15.515	2278	2283	2285	rBV	18614	23054	17.72%	2.498%
28	15.974	2355	2361	2368	rVB	4420	9003	6.92%	0.975%
29	16.404	2430	2434	2436	rBB2	1135	1527	1.17%	0.165%
30	16.486	2443	2448	2453	rBV3	2559	4147	3.19%	0.449%
31	17.080	2545	2549	2556	rBV2	1269	3043	2.34%	0.330%
32	18.645	2811	2815	2817	rBV	1132	1582	1.22%	0.171%

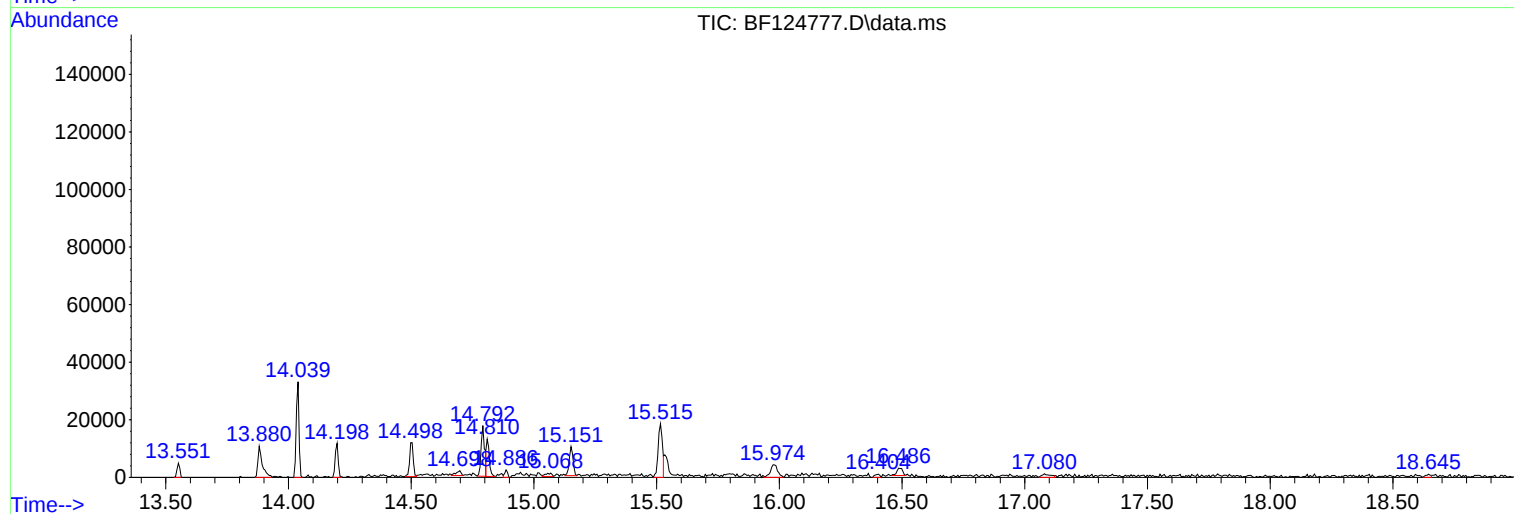
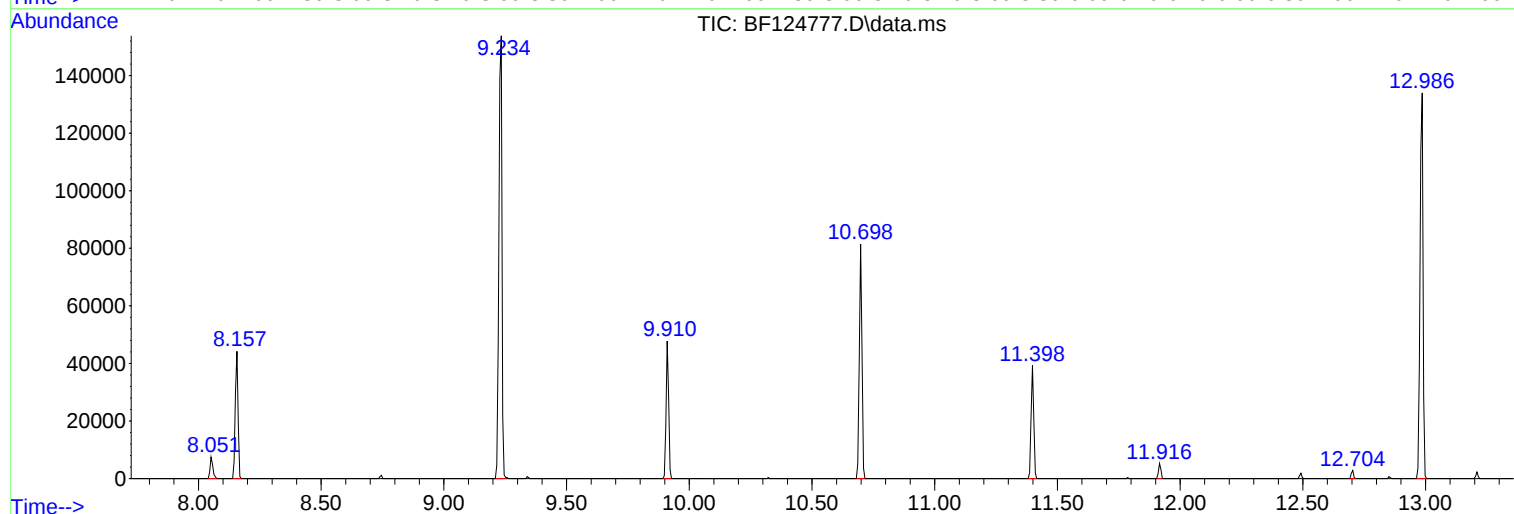
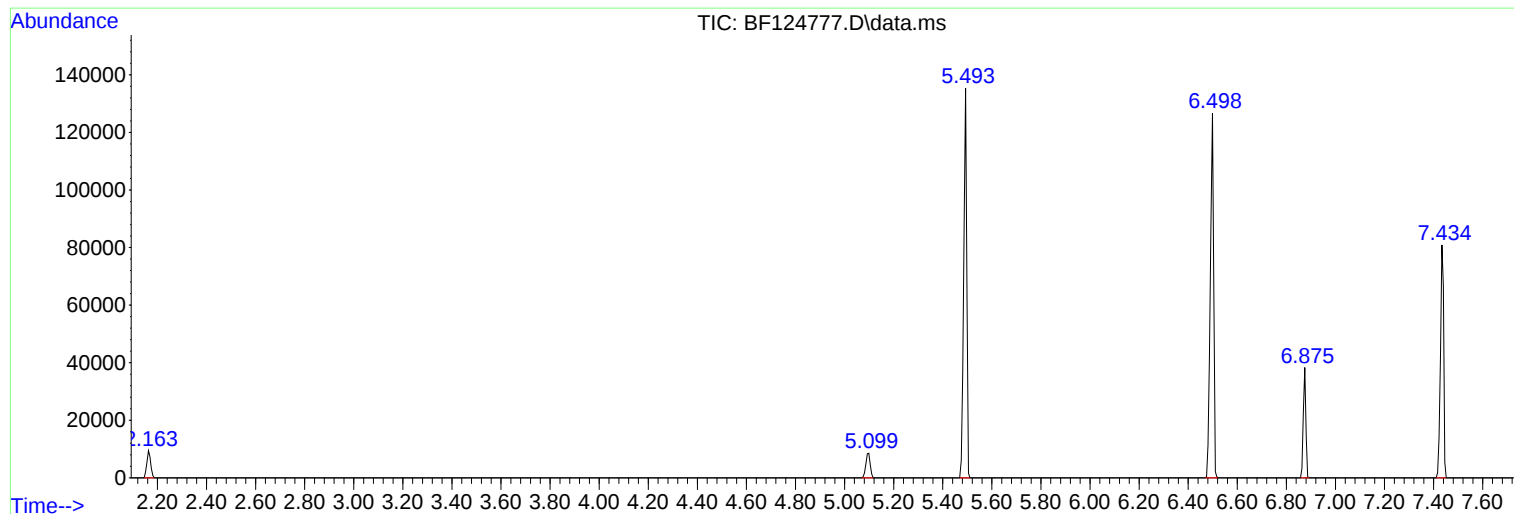
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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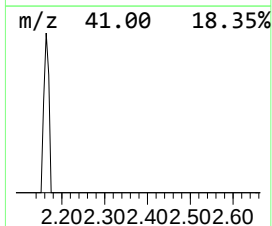
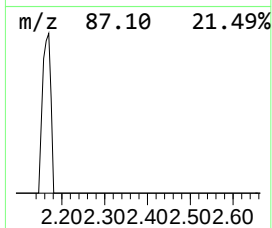
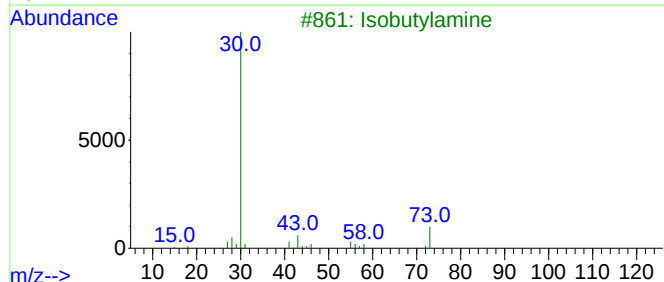
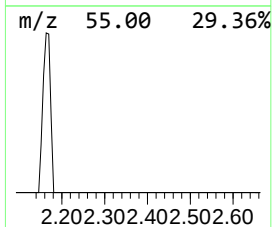
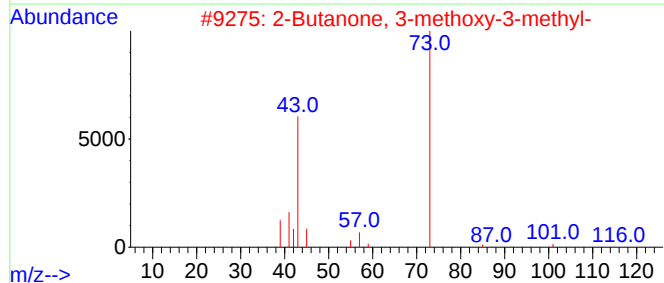
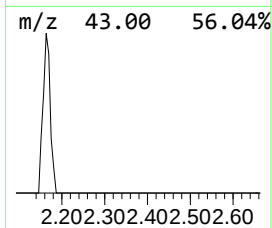
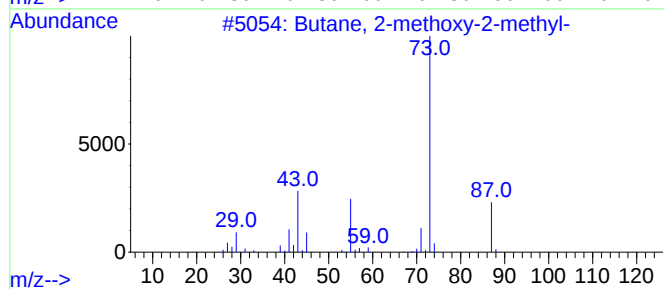
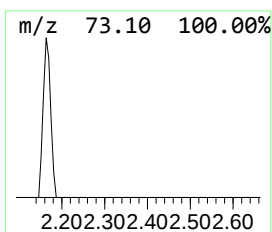
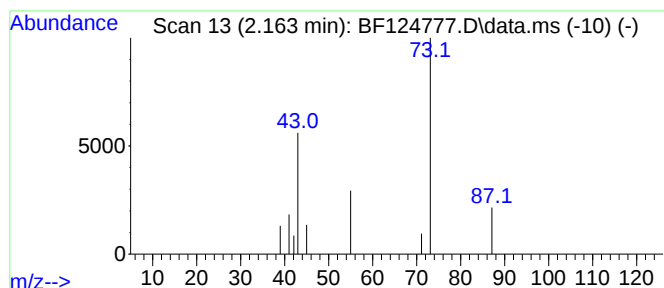
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.163	7.31 ng	10190	1,4-Dichlorobenzene-d4	6.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
2			2-Butanone, 3-methoxy-3-methyl-	116	C6H12O2	036687-98-6	9
3			Isobutylamine	73	C4H11N	000078-81-9	5
4			1-Pentanamine	87	C5H13N	000110-58-7	4
5			3,4-Hexanediol, 2,5-dimethyl-	146	C8H18O2	022607-11-0	4



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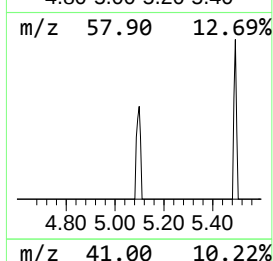
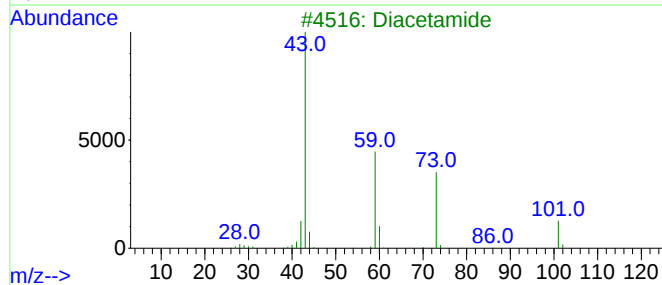
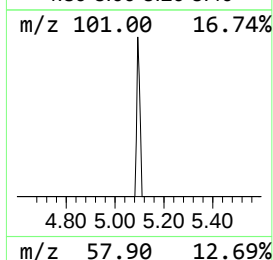
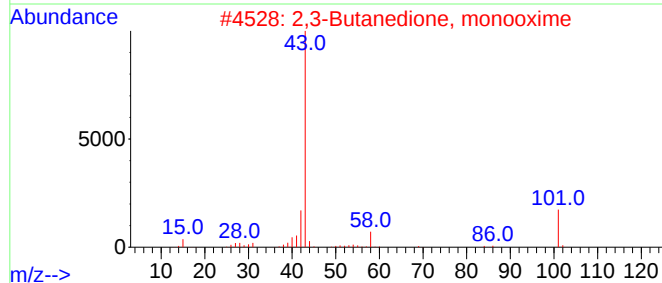
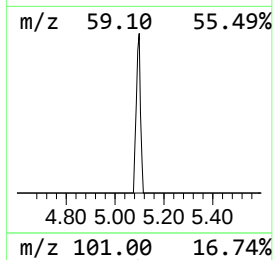
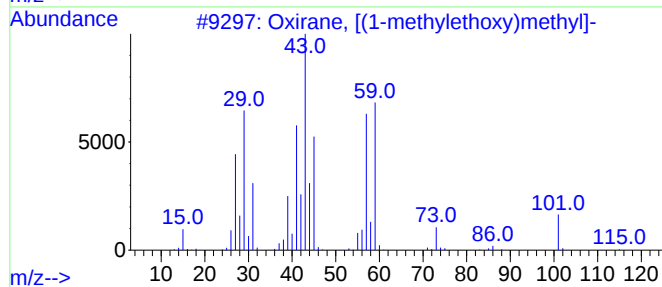
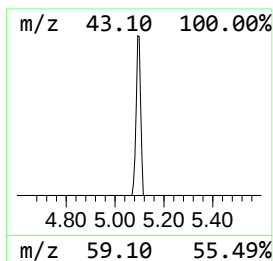
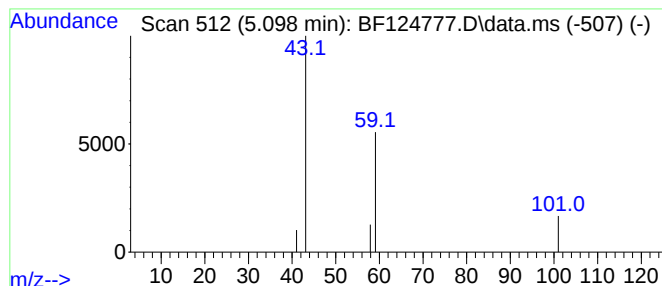
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown5.098 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.098	7.59 ng	10590	1,4-Dichlorobenzene-d4	6.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	9
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	7
3			Diacetamide	101	C4H7NO2	000625-77-4	7
4			Formamide, N-butyl-	101	C5H11NO	000871-71-6	4
5			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	4



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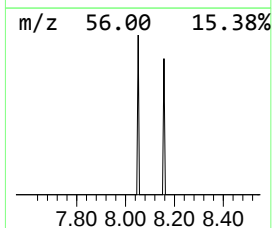
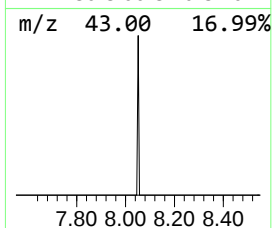
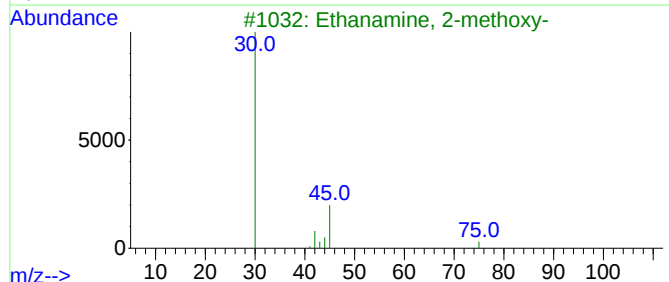
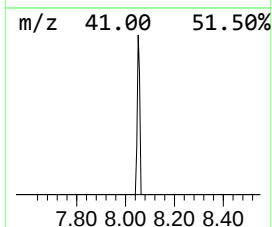
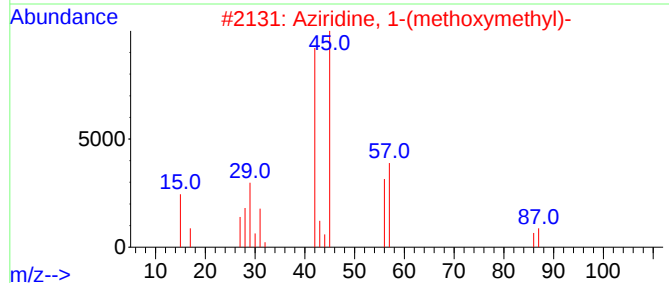
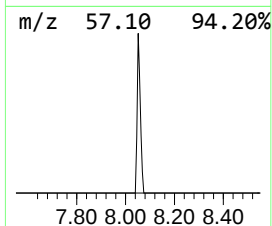
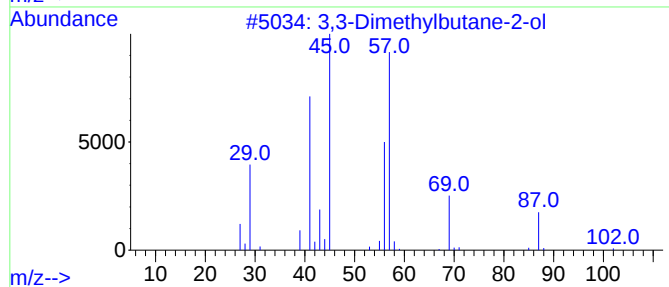
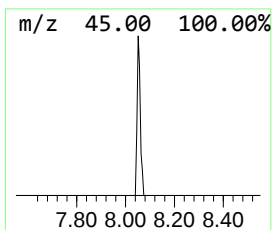
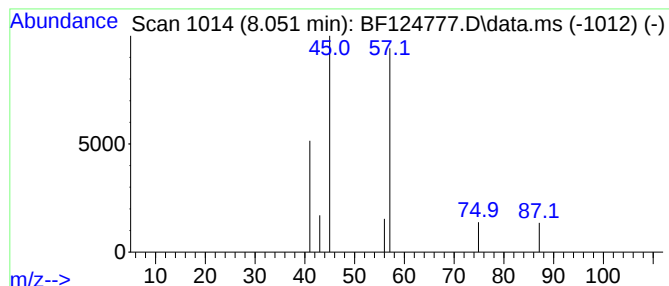
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown8.051 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.051	3.39 ng	5910	Naphthalene-d8	8.157

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	9
2			Aziridine, 1-(methoxymethyl)-	87	C4H9NO	001497-83-2	5
3			Ethanamine, 2-methoxy-	75	C3H9NO	000109-85-3	4
4			2-(Isobutoxymethyl)oxirane	130	C7H14O2	003814-55-9	4
5			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	4



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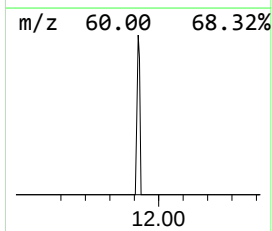
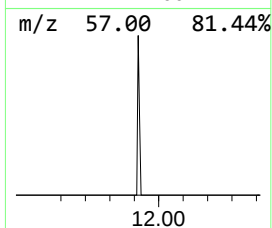
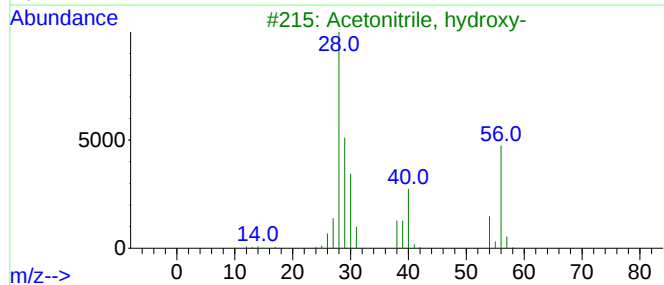
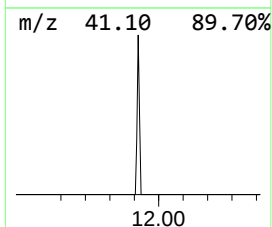
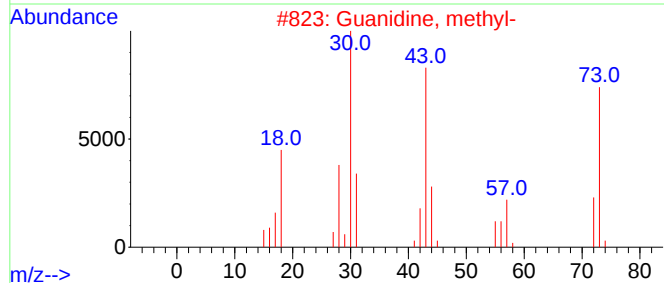
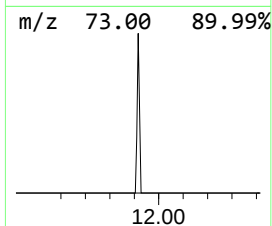
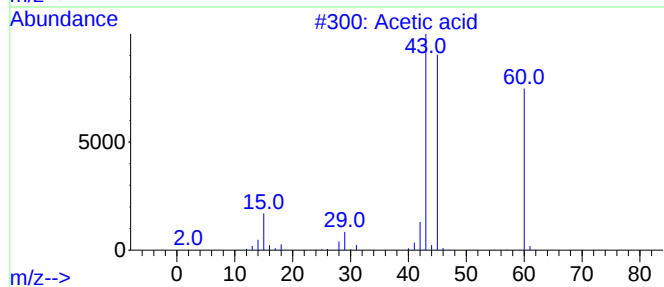
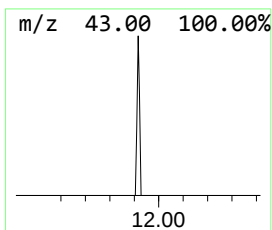
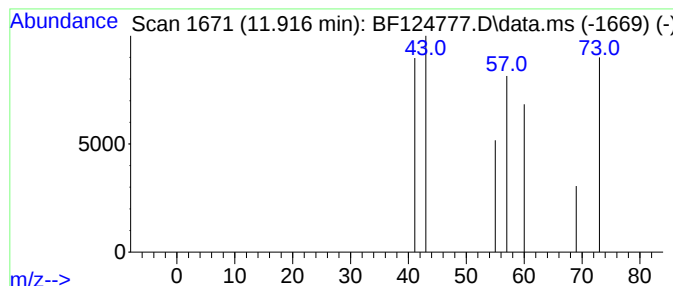
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 unknown11.916 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.916	2.43 ng	3586	Phenanthrene-d10		11.398	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Acetic acid		60	C2H4O2	000064-19-7	4
2	Guanidine, methyl-		73	C2H7N3	000471-29-4	4
3	Acetonitrile, hydroxy-		57	C2H3NO	000107-16-4	3
4	1-Propanol		60	C3H8O	000071-23-8	3
5	Isobutylamine		73	C4H11N	000078-81-9	3



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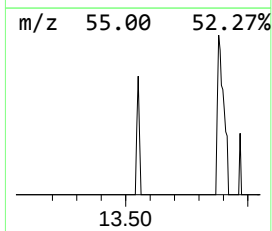
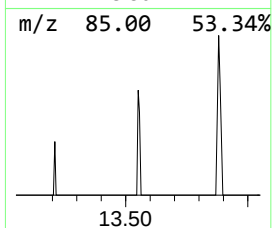
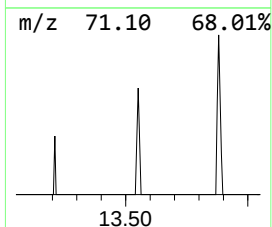
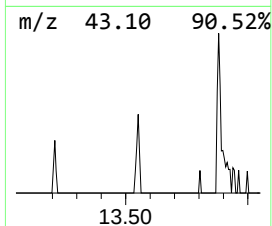
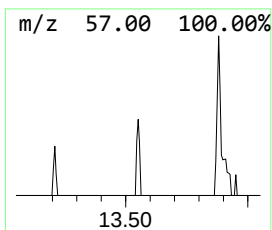
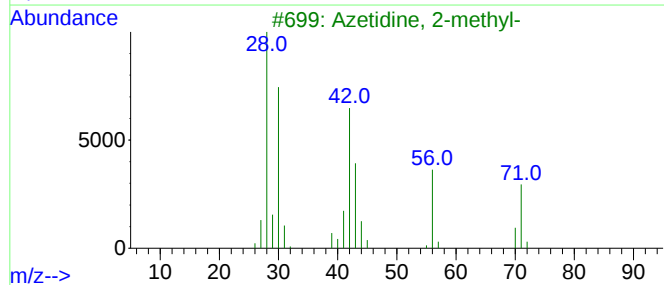
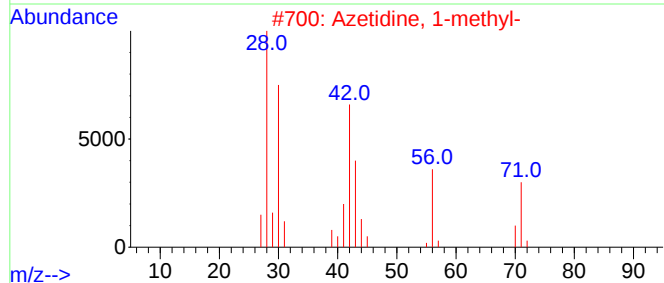
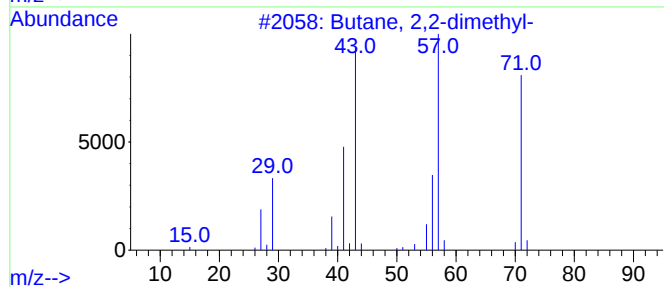
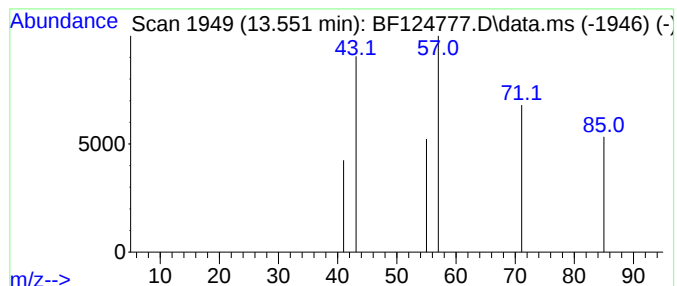
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TIC Library : C:\Database\NIST20.L
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Peak Number 5 unknown13.551 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.551	3.03 ng	3921	Chrysene-d12	14.039

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	9
2			Azetidine, 1-methyl-	71	C4H9N	004923-79-9	5
3			Azetidine, 2-methyl-	71	C4H9N	019812-49-8	5
4			1H-Tetrazol-5-amine	85	CH3N5	004418-61-5	4
5			5H-Tetrazol-5-amine	85	CH3N5	1000273-02-0	4



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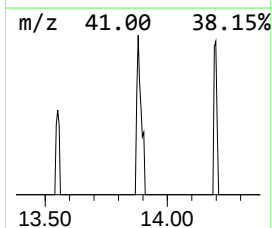
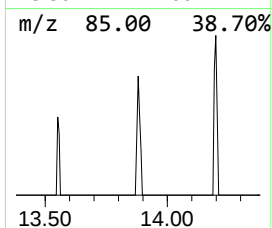
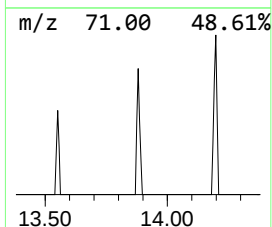
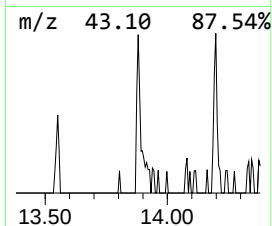
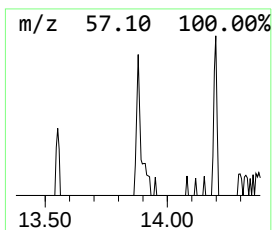
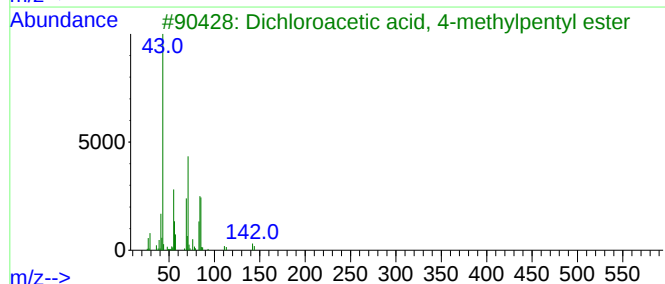
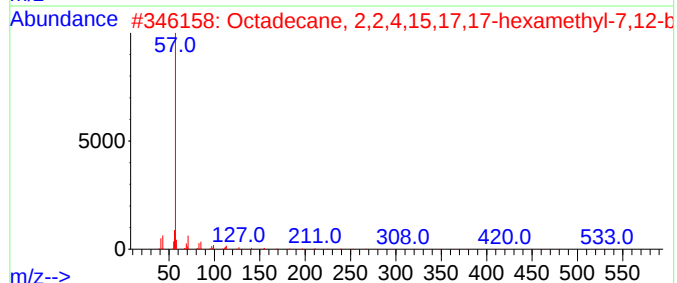
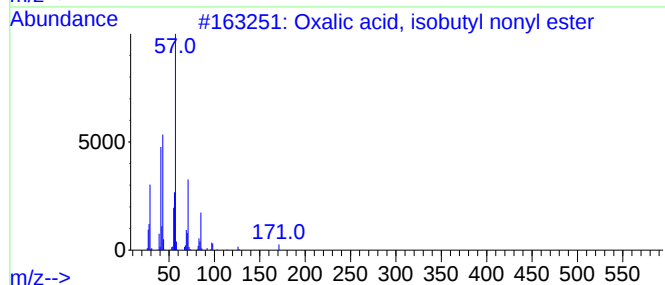
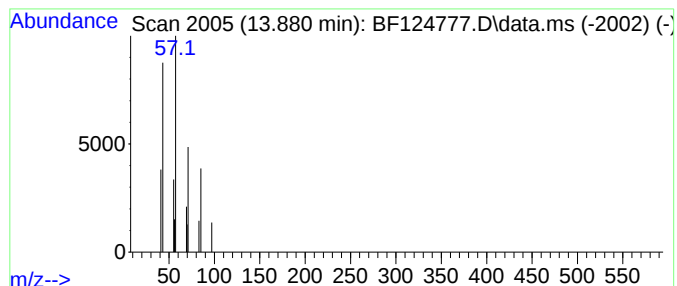
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown13.880 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.880	10.57 ng	13660	Chrysene-d12	14.039

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, isobutyl nonyl ester	272	C15H28O4	1010309-37-4	38
2			Octadecane, 2,2,4,15,17,17-hexam...	591	C42H86	055470-97-8	36
3			Dichloroacetic acid, 4-methylpen...	212	C8H14Cl2O2	1000282-43-7	25
4			1-Pentanol, 4-methyl-2-methylene-	114	C7H14O	121505-31-5	25
5			1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072621\
Data File : BF124777.D
Acq On : 26 Jul 2021 22:05
Operator : JU/CG
Sample : M3151-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMP-1

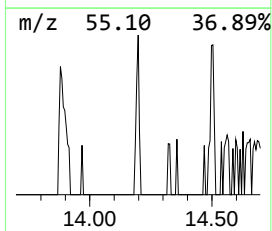
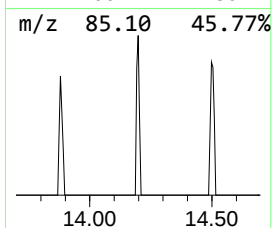
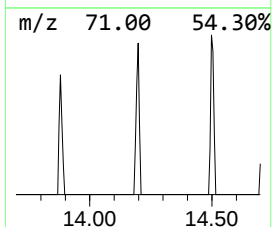
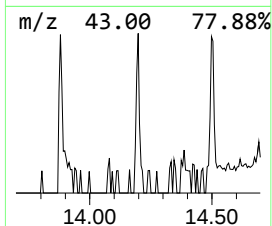
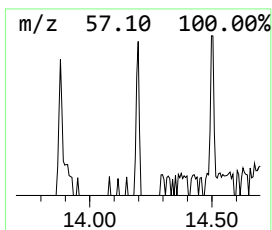
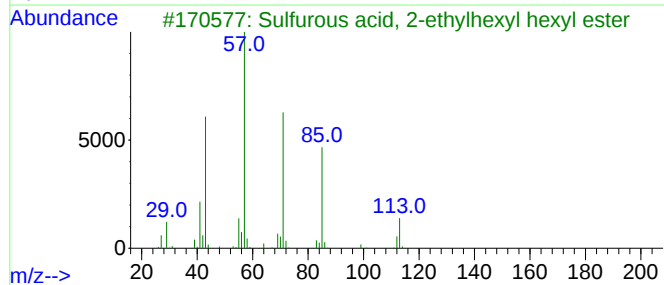
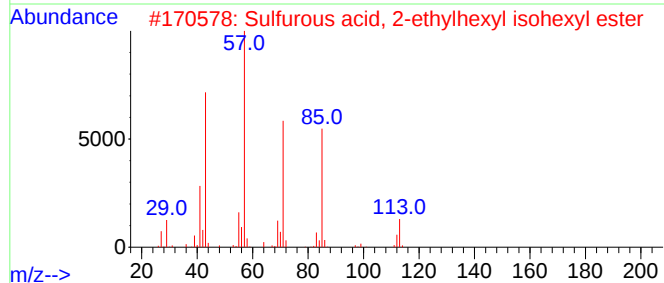
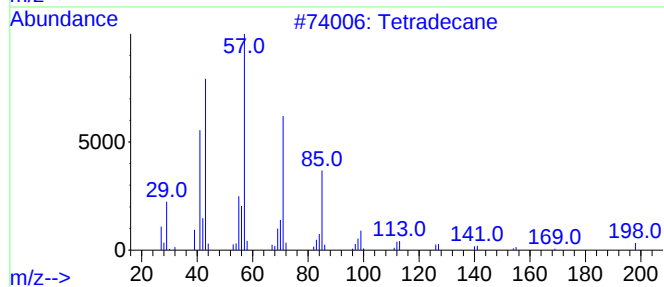
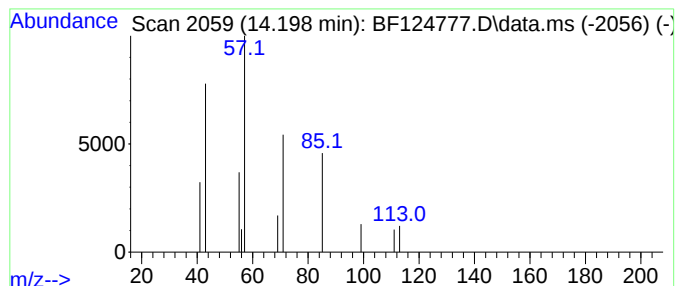
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tetradecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.198	7.77 ng	10048	Chrysene-d12	14.039

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	78
2			Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	72
3			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	64
4			Octane, 2,3,3-trimethyl-	156	C11H24	062016-30-2	50
5			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072621\
Data File : BF124777.D
Acq On : 26 Jul 2021 22:05
Operator : JU/CG
Sample : M3151-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMP-1

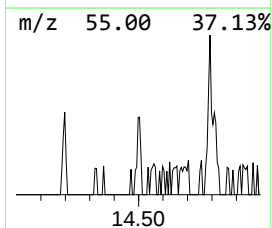
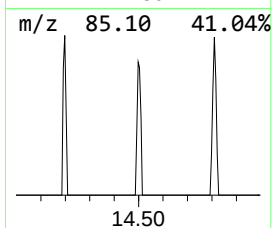
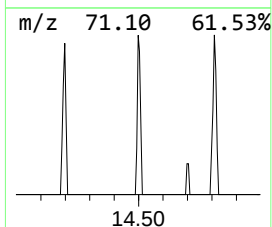
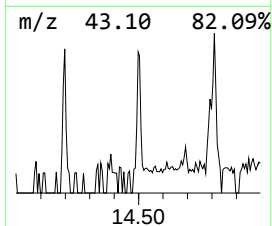
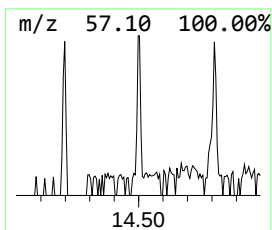
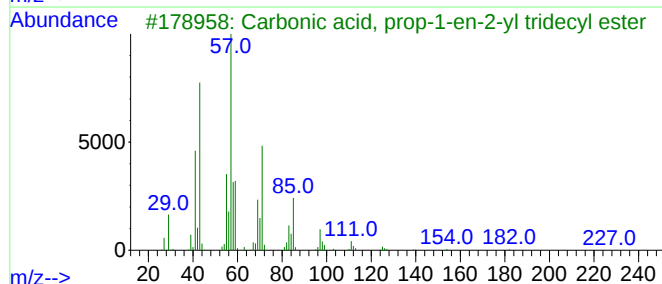
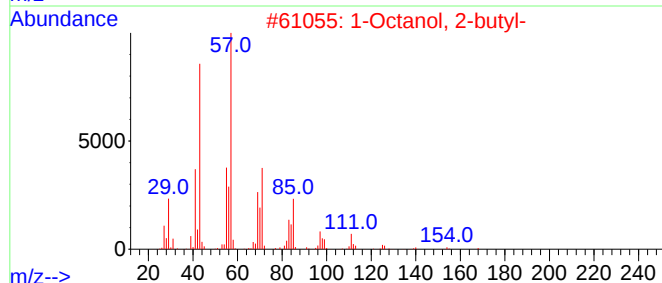
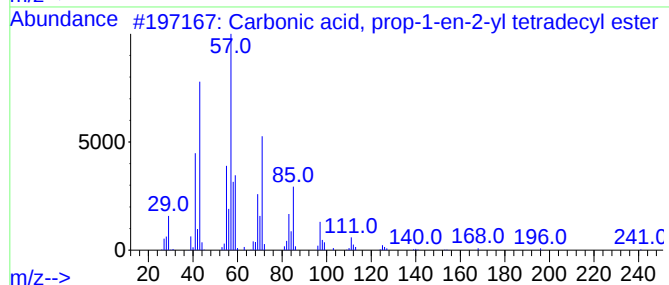
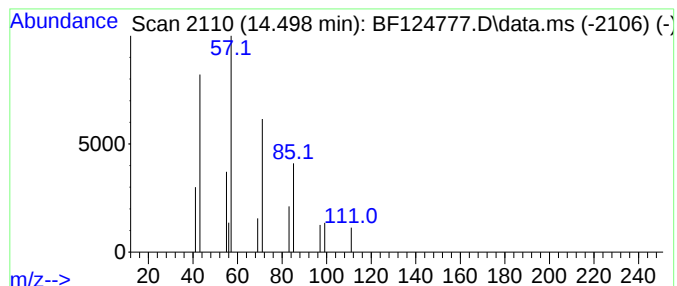
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Carbonic acid, prop-1-en-2-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.498	9.39 ng	12136	Chrysene-d12	14.039

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	78
2			1-Octanol, 2-butyl-	186	C12H26O	003913-02-8	39
3			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	38
4			1-Heptadecanamine	255	C17H37N	004200-95-7	38
5			Pentane, 3-ethyl-2,3-dimethyl-	128	C9H20	016747-33-4	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072621\
Data File : BF124777.D
Acq On : 26 Jul 2021 22:05
Operator : JU/CG
Sample : M3151-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

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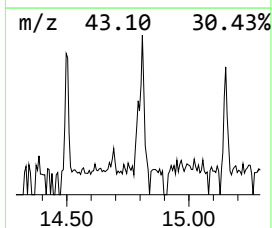
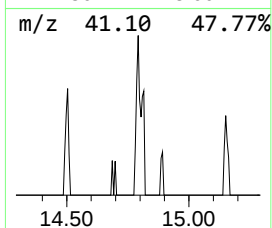
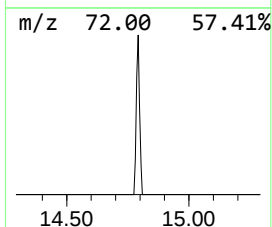
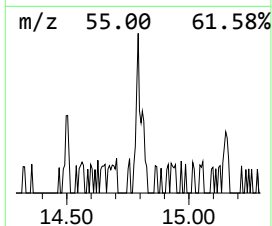
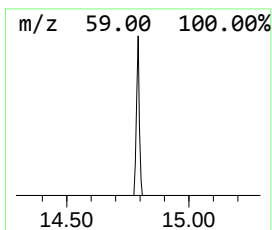
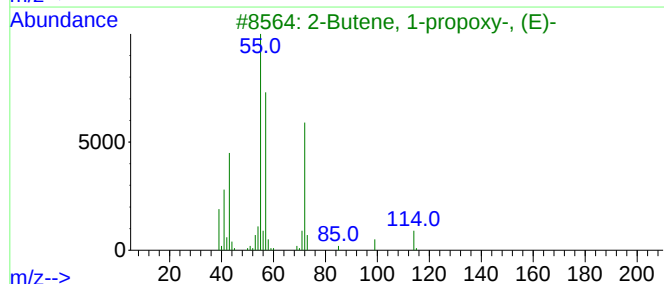
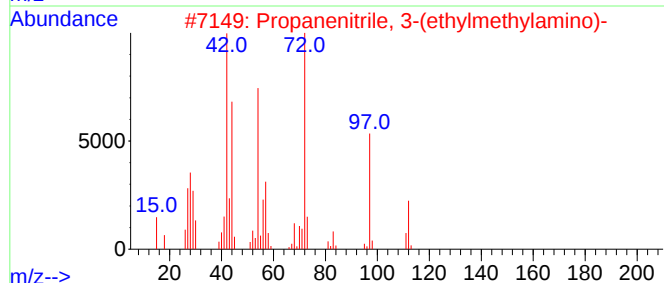
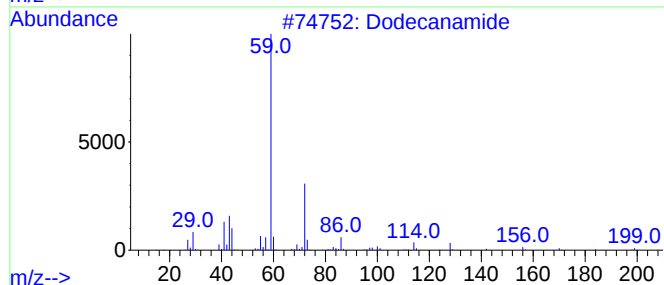
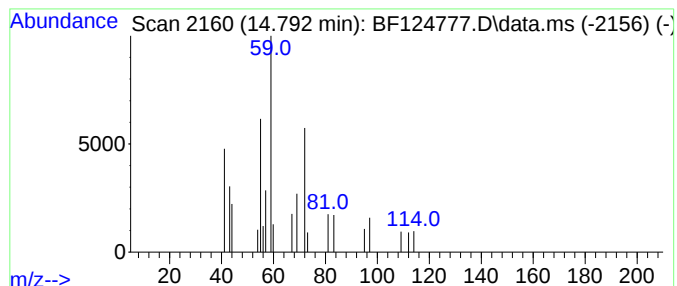
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown14.792 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.792	15.74 ng	18149	Perylene-d12	15.515

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecanamide	199	C12H25NO	001120-16-7	38
2			Propanenitrile, 3-(ethylmethy...)	112	C6H12N2	055619-09-5	10
3			2-Butene, 1-propoxy-, (E)-	114	C7H14O	056052-71-2	10
4			2-Propenoic acid	72	C3H4O2	000079-10-7	9
5			Cyclopentylsilane	100	C5H12Si	080249-74-7	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072621\
Data File : BF124777.D
Acq On : 26 Jul 2021 22:05
Operator : JU/CG
Sample : M3151-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

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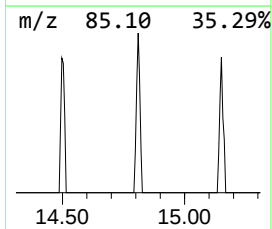
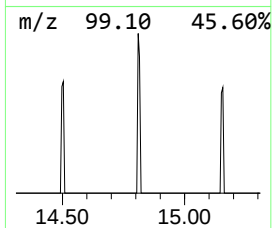
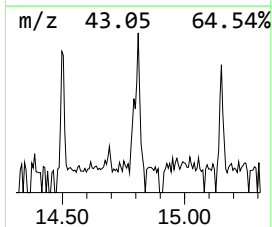
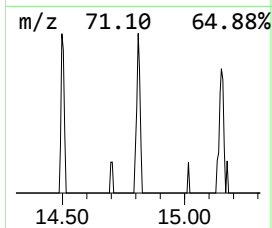
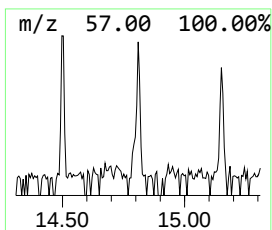
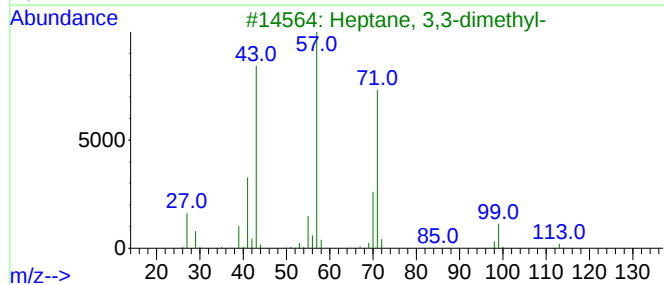
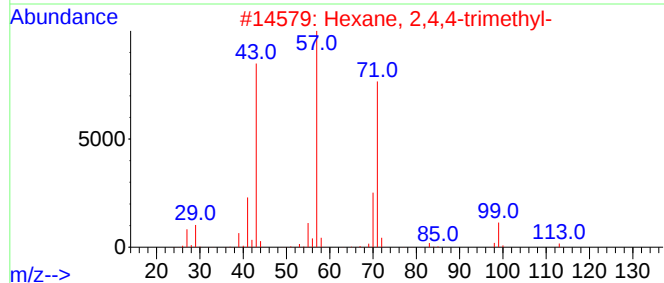
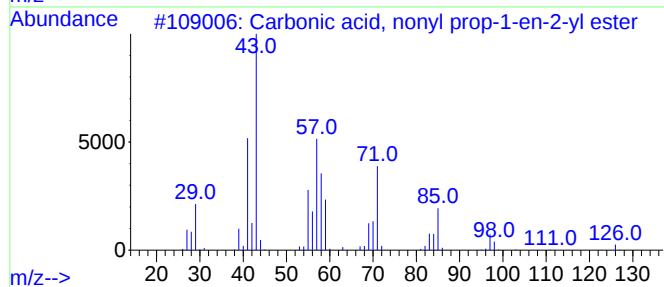
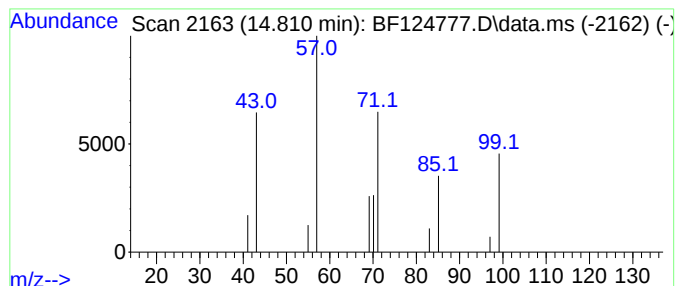
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown14.810 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.810	8.45 ng	9742	Perylene-d12	15.515

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, nonyl prop-1-en-2...	228	C13H24O3	1000382-53-9	36
2			Hexane, 2,4,4-trimethyl-	128	C9H20	016747-30-1	33
3			Heptane, 3,3-dimethyl-	128	C9H20	004032-86-4	25
4			2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	9
5			Hexane, 1-(hexyloxy)-5-methyl-	200	C13H28O	074421-19-5	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072621\
Data File : BF124777.D
Acq On : 26 Jul 2021 22:05
Operator : JU/CG
Sample : M3151-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
COMP-1

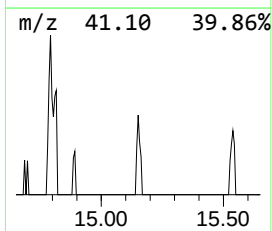
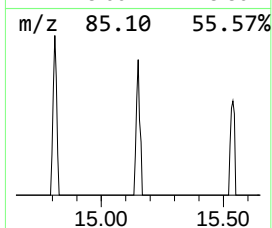
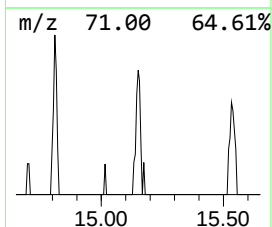
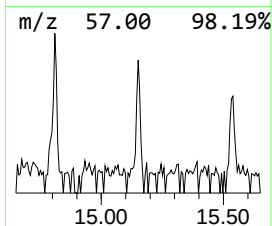
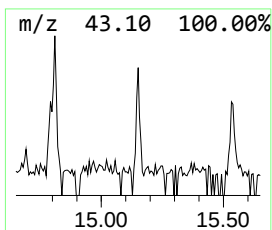
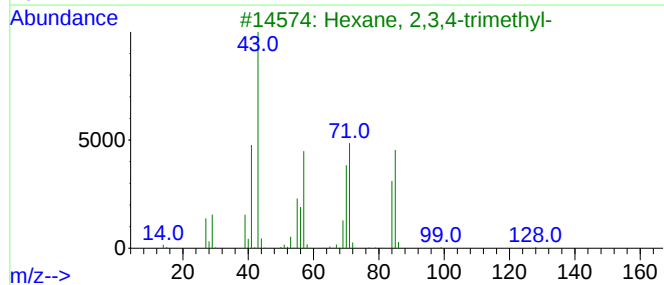
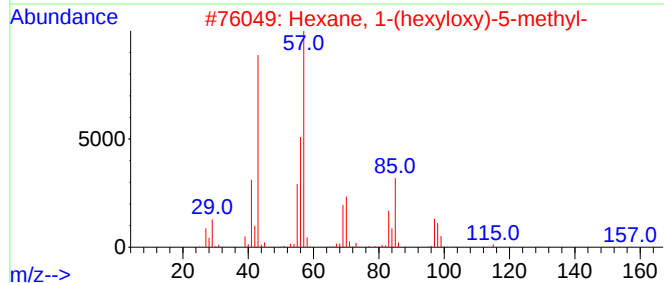
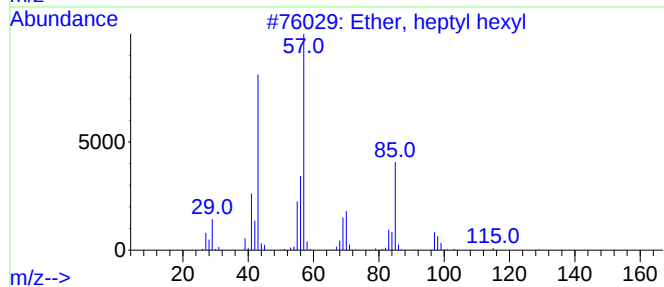
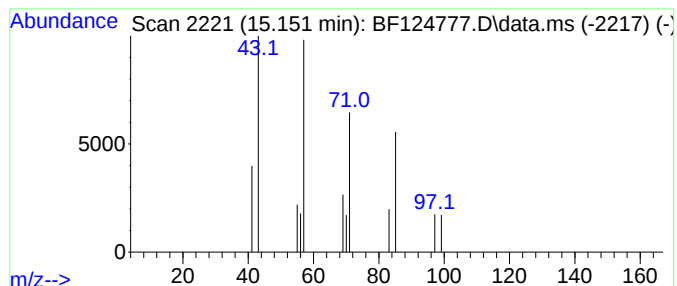
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown15.151 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.151	9.35 ng	10774	Perylene-d12	15.515

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ether, heptyl hexyl	200	C13H28O	007289-40-9	47
2			Hexane, 1-(hexyloxy)-5-methyl-	200	C13H28O	074421-19-5	38
3			Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	28
4			Dichloroacetic acid, 4-methylpen...	212	C8H14Cl2O2	1000282-43-7	28
5			Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	28



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072621\
Data File : BF124777.D
Acq On : 26 Jul 2021 22:05
Operator : JU/CG
Sample : M3151-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

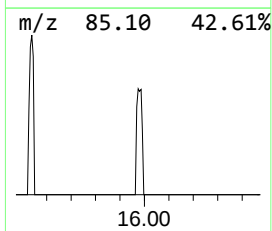
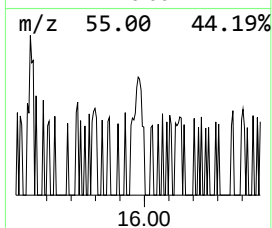
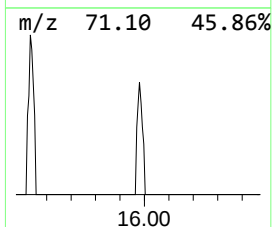
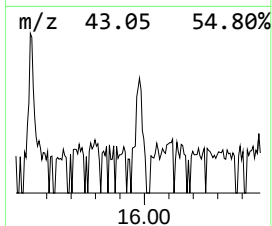
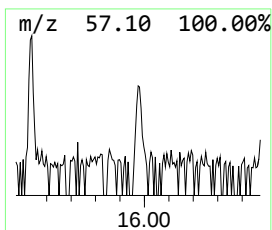
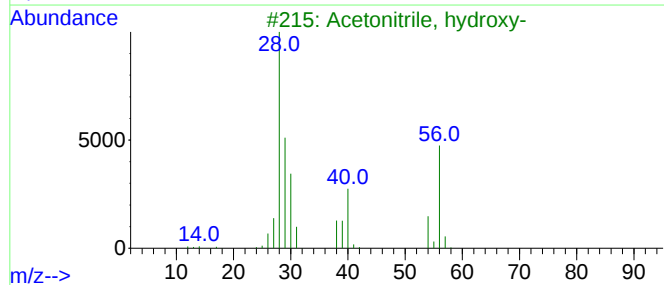
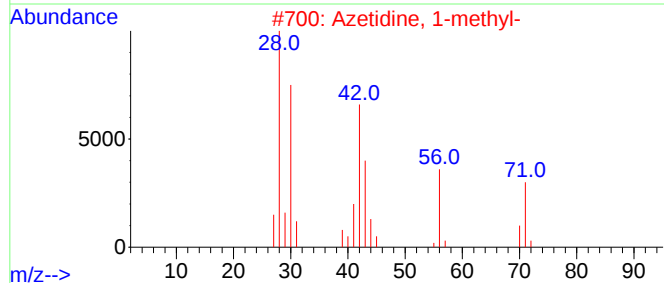
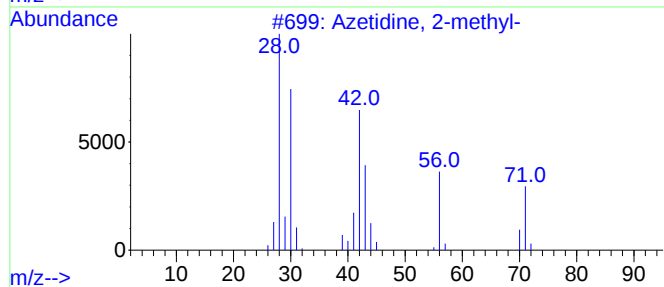
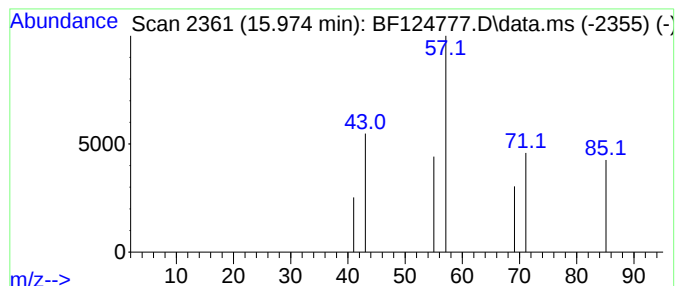
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown15.974 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.974	7.81 ng	9003	Perylene-d12		15.515	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Azetidine, 2-methyl-		71	C4H9N	019812-49-8	5
2	Azetidine, 1-methyl-		71	C4H9N	004923-79-9	5
3	Acetonitrile, hydroxy-		57	C2H3NO	000107-16-4	4
4	5H-Tetrazol-5-amine		85	CH3N5	1000273-02-0	4
5	1H-Tetrazol-5-amine		85	CH3N5	004418-61-5	4



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072621\
Data File : BF124777.D
Acq On : 26 Jul 2021 22:05
Operator : JU/CG
Sample : M3151-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMP-1

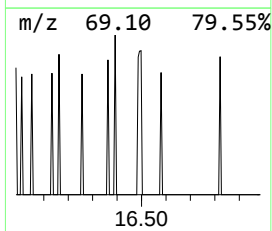
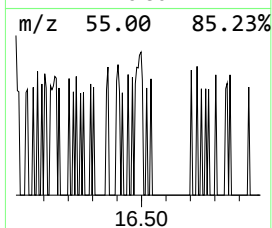
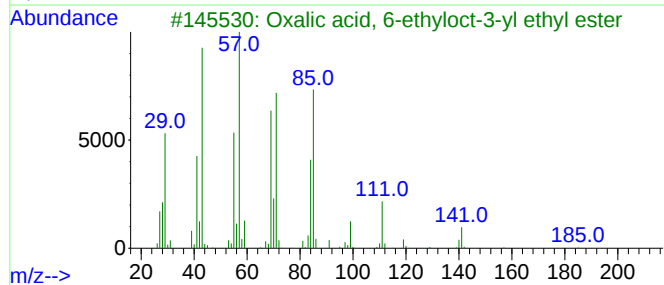
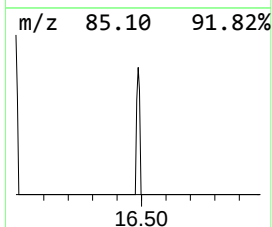
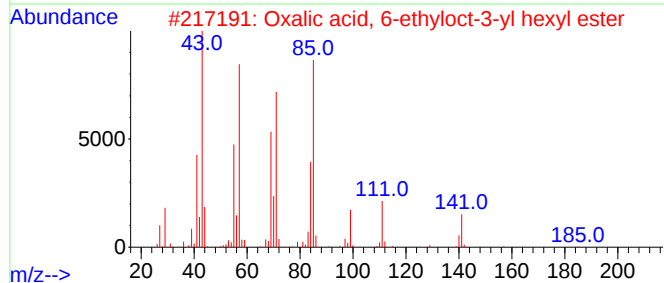
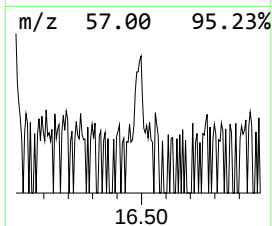
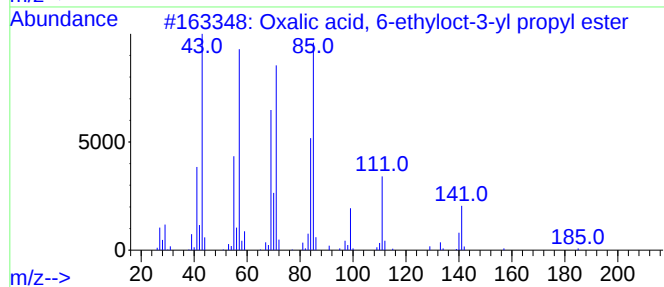
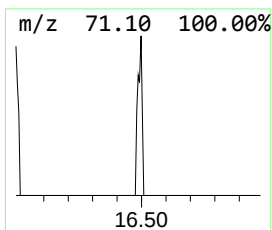
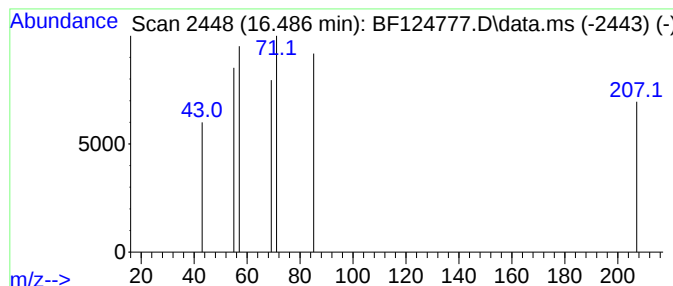
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown16.486 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.486	3.60 ng	4147	Perylene-d12	15.515

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, 6-ethyloct-3-yl pro...	272	C15H28O4	1000309-34-0	40
2			Oxalic acid, 6-ethyloct-3-yl hex...	314	C18H34O4	1000309-34-4	9
3			Oxalic acid, 6-ethyloct-3-yl eth...	258	C14H26O4	1000309-33-9	9
4			Oxalic acid, 6-ethyloct-3-yl iso...	314	C18H34O4	1000309-34-3	9
5			Oxalic acid, 6-ethyloct-3-yl hep...	328	C19H36O4	1000309-34-5	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072621\
Data File : BF124777.D
Acq On : 26 Jul 2021 22:05
Operator : JU/CG
Sample : M3151-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMP-1

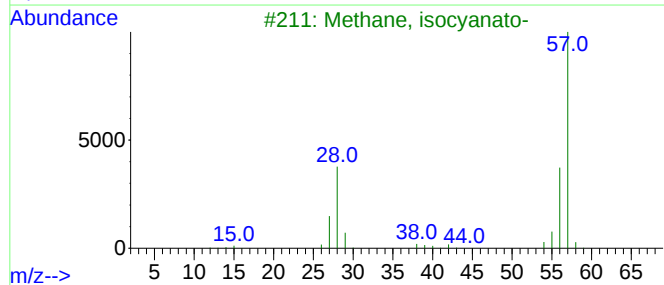
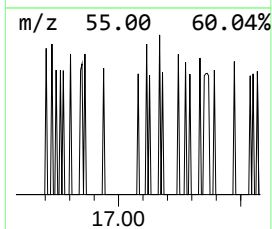
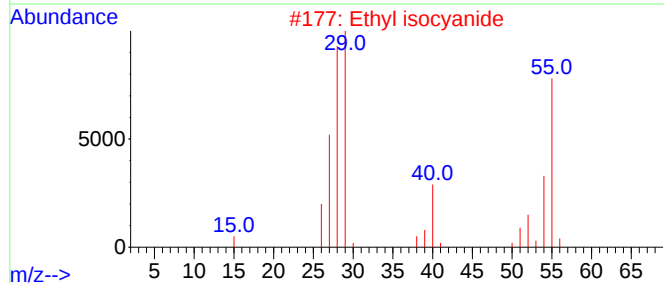
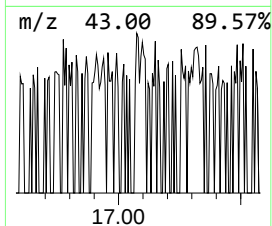
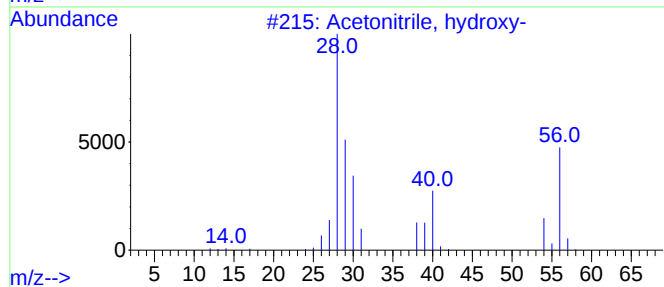
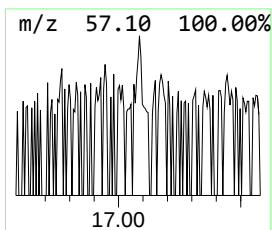
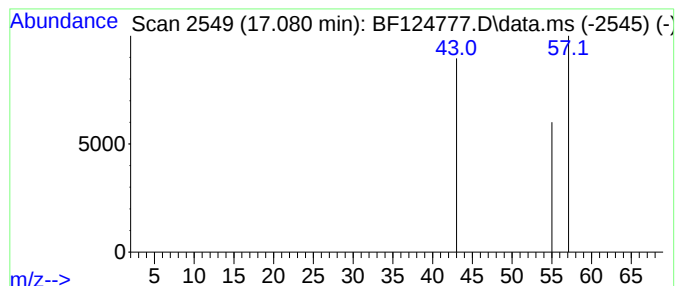
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown17.080 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.080	2.64 ng	3043	Perylene-d12	15.515

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Acetonitrile, hydroxy-	57	C2H3NO	000107-16-4	3
2		Ethyl isocyanide	55	C3H5N	000624-79-3	1
3		Methane, isocyanato-	57	C2H3NO	000624-83-9	1
4		Azetidine	57	C3H7N	000503-29-7	1
5		2-Propen-1-amine	57	C3H7N	000107-11-9	1



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072621\
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 Operator : JU/CG
 Sample : M3151-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072221.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
Butane, 2-metho...	2.163	7.3	ng	10190	1	6.875	27893	20.0
unknown5.098	5.098	7.6	ng	10590	1	6.875	27893	20.0
unknown8.051	8.051	3.4	ng	5910	2	8.157	34882	20.0
unknown11.916	11.916	2.4	ng	3586	4	11.398	29516	20.0
unknown13.551	13.551	3.0	ng	3921	5	14.039	25847	20.0
unknown13.880	13.880	10.6	ng	13660	5	14.039	25847	20.0
Tetradecane	14.198	7.8	ng	10048	5	14.039	25847	20.0
Carbonic acid, ...	14.498	9.4	ng	12136	5	14.039	25847	20.0
unknown14.792	14.792	15.7	ng	18149	6	15.515	23054	20.0
unknown14.810	14.810	8.4	ng	9742	6	15.515	23054	20.0
unknown15.151	15.151	9.3	ng	10774	6	15.515	23054	20.0
unknown15.974	15.974	7.8	ng	9003	6	15.515	23054	20.0
unknown16.486	16.486	3.6	ng	4147	6	15.515	23054	20.0
unknown17.080	17.080	2.6	ng	3043	6	15.515	23054	20.0