

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042722\
 Data File : BF127956.D
 Acq On : 27 Apr 2022 20:24
 Operator : CG\JU
 Sample : N2482-15
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SPN-8

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: BF127956.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	330	334	339	rBV	48309	59526	1.69%	0.307%
2	4.457	365	369	377	rBV	347360	403347	11.44%	2.080%
3	5.157	484	488	495	rBV	121135	142487	4.04%	0.735%
4	5.422	530	533	537	rVB	51905	45979	1.30%	0.237%
5	5.545	550	554	558	rBV	1636298	1465569	41.58%	7.559%
6	6.104	645	649	652	rVB	70808	65013	1.84%	0.335%
7	6.157	654	658	662	rVB	73025	59606	1.69%	0.307%
8	6.387	694	697	700	rBV	128183	114919	3.26%	0.593%
9	6.487	711	714	716	rBV	50681	41532	1.18%	0.214%
10	6.545	720	724	733	rVV	1048465	947296	26.88%	4.886%
11	6.610	733	735	738	rVV	84566	73704	2.09%	0.380%
12	6.728	751	755	758	rBV	144979	172934	4.91%	0.892%
13	6.757	758	760	770	rVV	151092	202186	5.74%	1.043%
14	6.851	773	776	786	rBV	172119	265614	7.54%	1.370%
15	6.928	786	789	793	rVB	736865	597250	16.95%	3.081%
16	6.981	796	798	801	rBV	77309	71526	2.03%	0.369%
17	7.057	801	811	812	rBV3	23273	57525	1.63%	0.297%
18	7.110	817	820	823	rVB	119219	101334	2.88%	0.523%
19	7.334	853	858	865	rVB3	42675	64377	1.83%	0.332%
20	7.463	876	880	881	rBV	76670	70021	1.99%	0.361%
21	7.492	881	885	888	rVB	2152672	2132334	60.50%	10.998%
22	7.569	895	898	902	rBV	40882	45561	1.29%	0.235%
23	7.945	957	962	965	rVB2	59639	58800	1.67%	0.303%
24	8.175	993	1001	1003	rBV5	25369	49936	1.42%	0.258%
25	8.210	1003	1007	1011	rBV	850410	773291	21.94%	3.988%
26	8.463	1039	1050	1053	rBV5	21015	66370	1.88%	0.342%
27	9.204	1172	1176	1178	rBV	38363	51163	1.45%	0.264%
28	9.286	1185	1190	1194	rBV	3454009	3524565	100.00%	18.179%
29	9.386	1204	1207	1217	rVB	492077	585356	16.61%	3.019%
30	9.969	1303	1306	1310	rVB	970371	816431	23.16%	4.211%
31	10.380	1372	1376	1379	rVB	145032	123518	3.50%	0.637%
32	10.675	1422	1426	1428	rBV	113967	106422	3.02%	0.549%
33	10.763	1437	1441	1444	rBV	1920597	1703498	48.33%	8.786%
34	11.463	1556	1560	1563	rBV	831284	707318	20.07%	3.648%
35	11.839	1621	1624	1627	rVB	87408	62982	1.79%	0.325%
36	13.051	1825	1830	1833	rBV	2645080	2395449	67.96%	12.355%

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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

37	14.110	2006	2010	2019	rVB	534189	520943	14.78%	2.687%
38	15.615	2261	2266	2273	rBV2	484382	642397	18.23%	3.313%

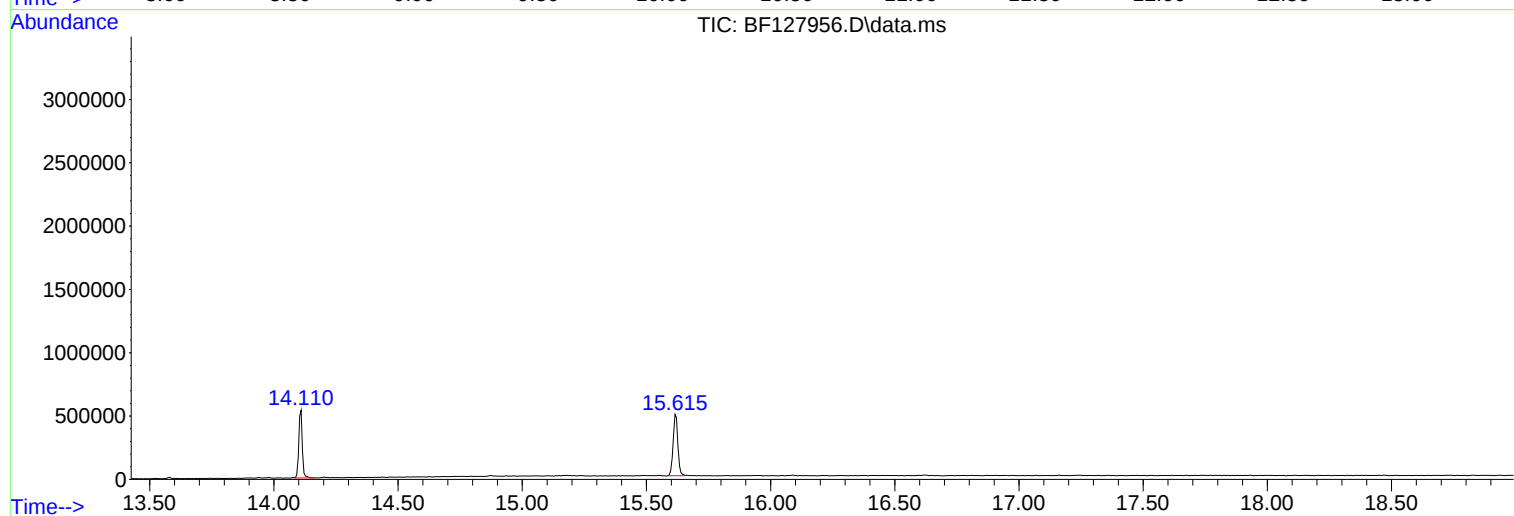
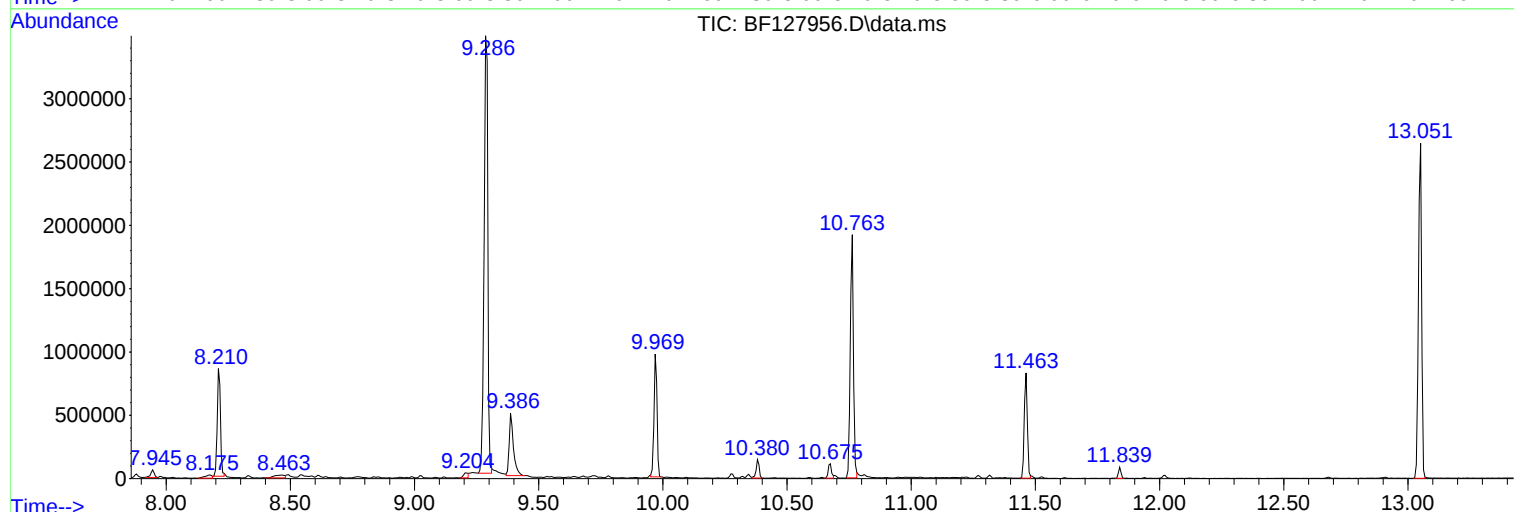
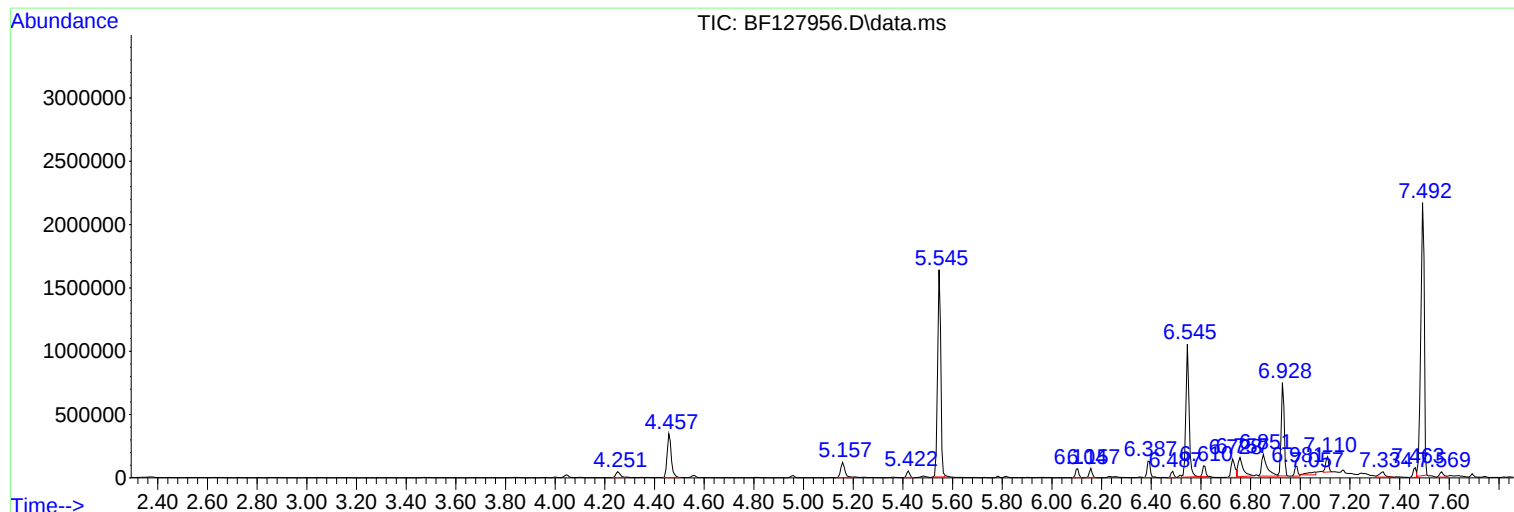
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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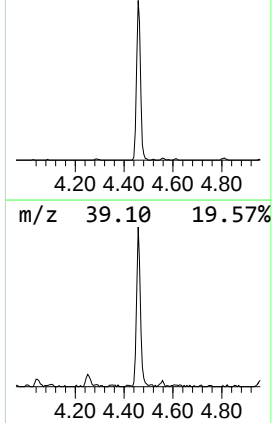
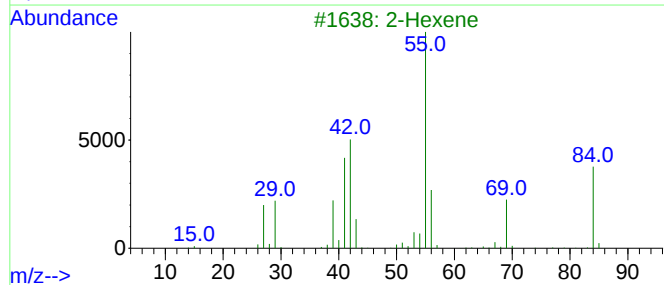
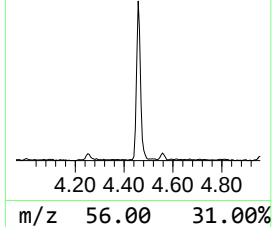
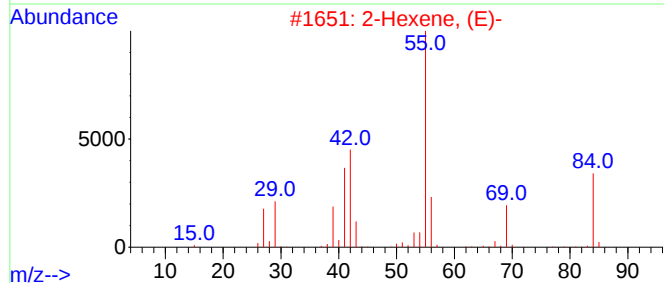
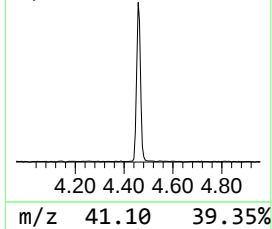
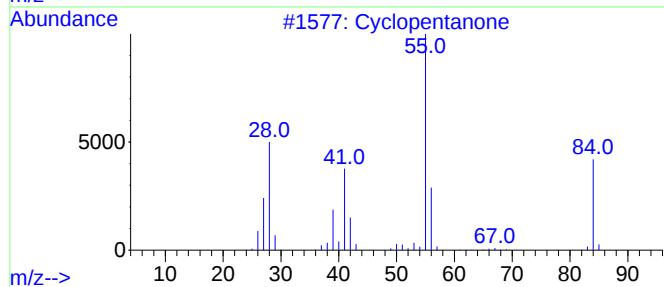
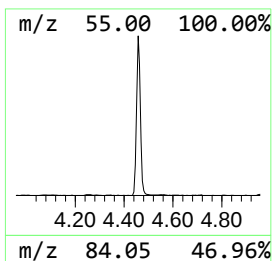
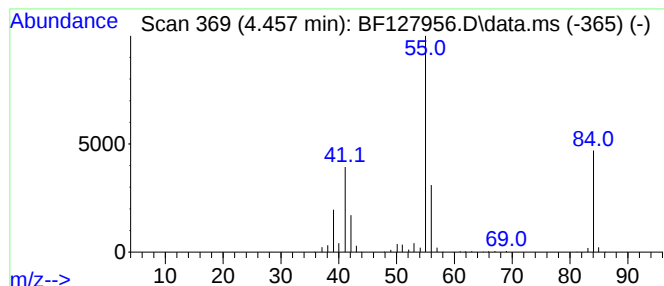
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Cyclopentanone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.457	13.51 ng	403347	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentanone	84	C5H8O	000120-92-3	90
2			2-Hexene, (E)-	84	C6H12	004050-45-7	64
3			2-Hexene	84	C6H12	000592-43-8	64
4			1H-Tetrazole, 1-methyl-	84	C2H4N4	016681-77-9	53
5			3-Hexene, (Z)-	84	C6H12	007642-09-3	50



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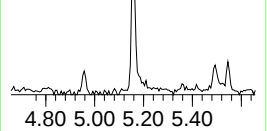
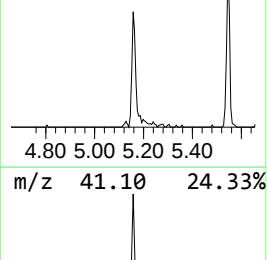
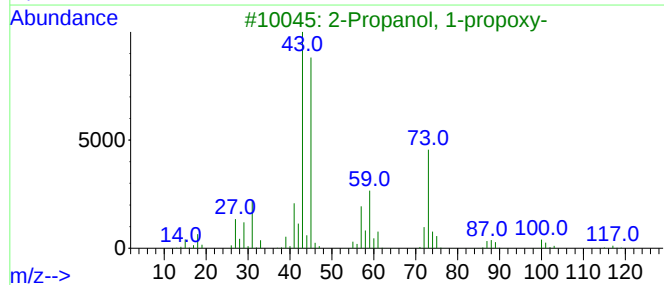
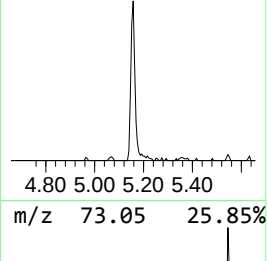
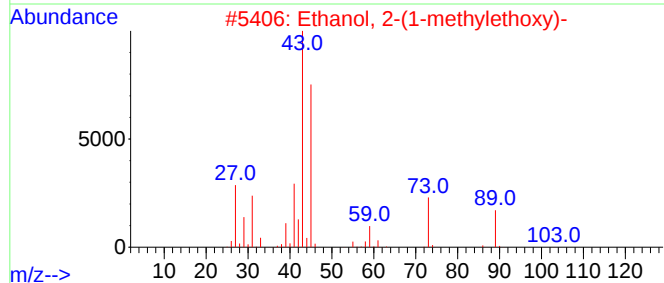
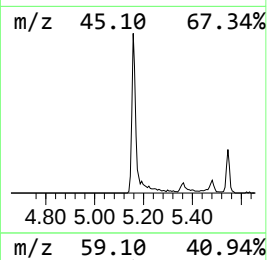
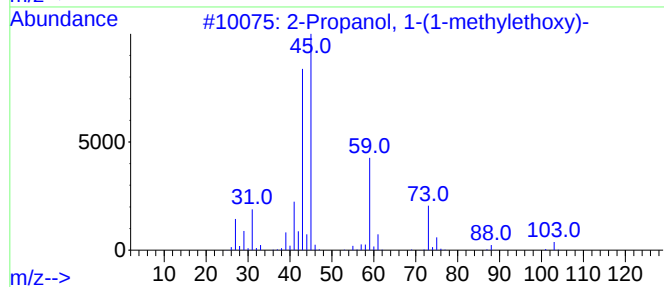
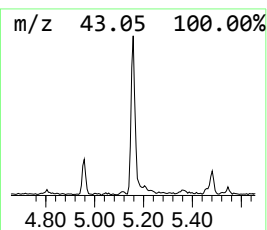
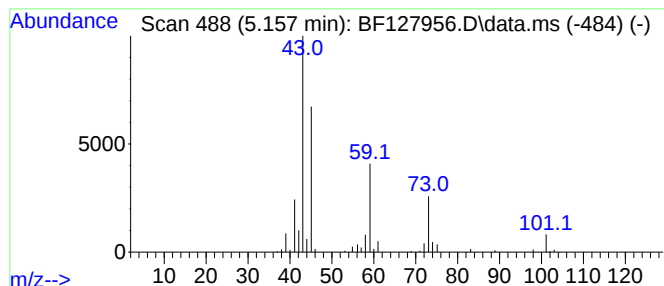
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042222.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.157	4.77 ng	142487	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	53		
2	Ethanol, 2-(1-methylethoxy)-	104	C5H12O2	000109-59-1	50		
3	2-Propanol, 1-propoxy-	118	C6H14O2	001569-01-3	40		
4	Butane, 1-(1-methylethoxy)-	116	C7H16O	001860-27-1	40		
5	Butanoic acid, 4-methoxy-	118	C5H10O3	029006-02-8	40		



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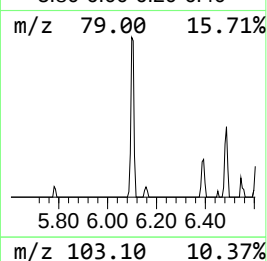
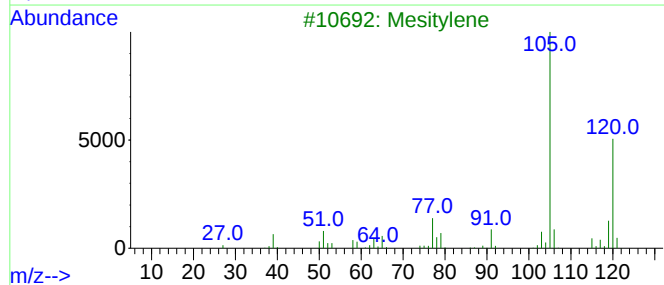
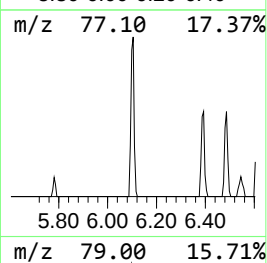
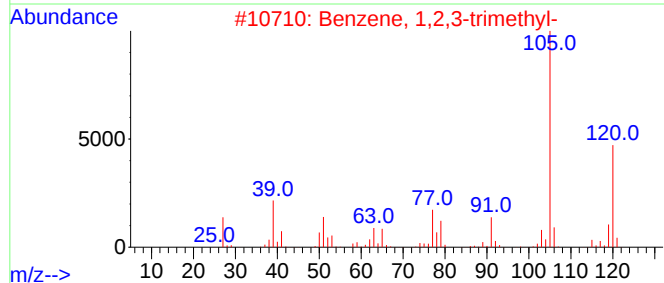
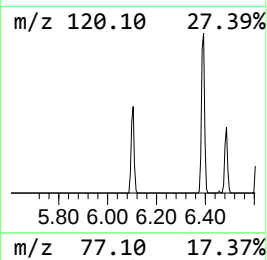
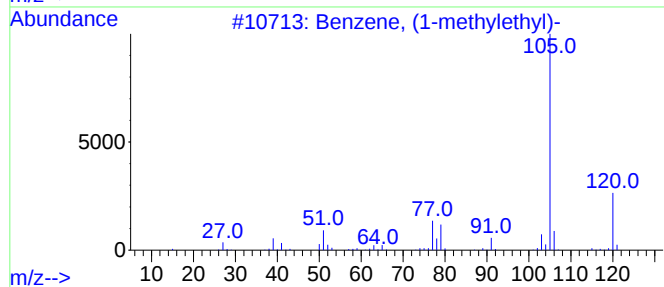
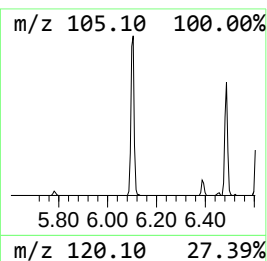
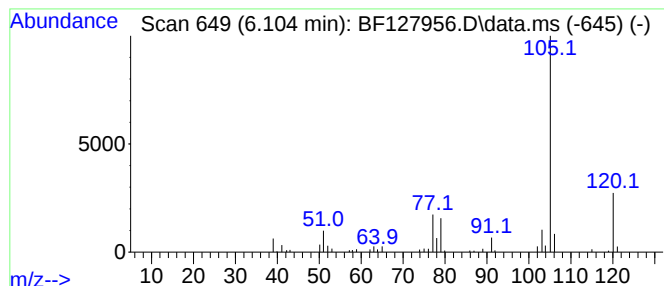
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Benzene, (1-methylethyl)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.104	2.18 ng	65013	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3			Mesitylene	120	C9H12	000108-67-8	86
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	83
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	80



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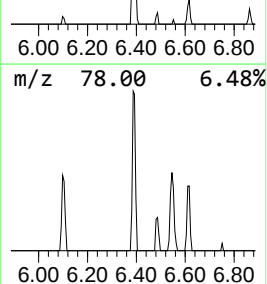
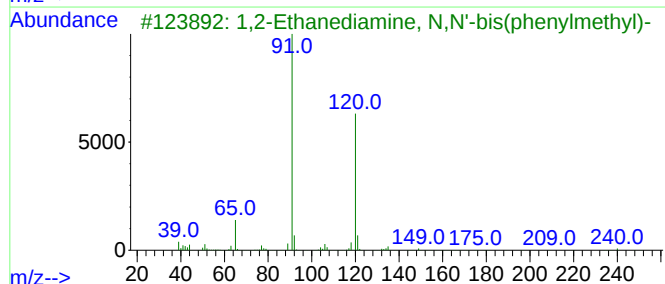
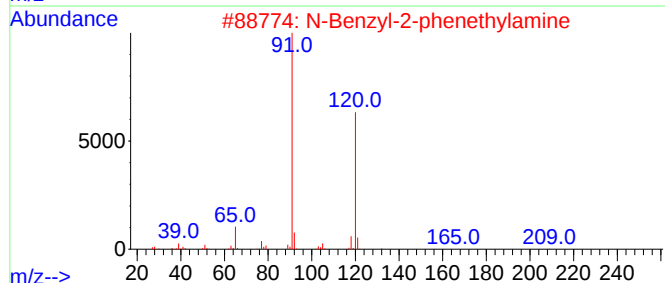
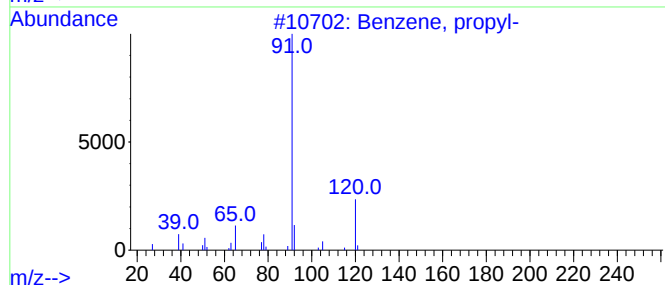
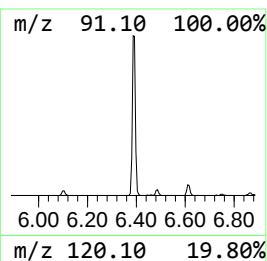
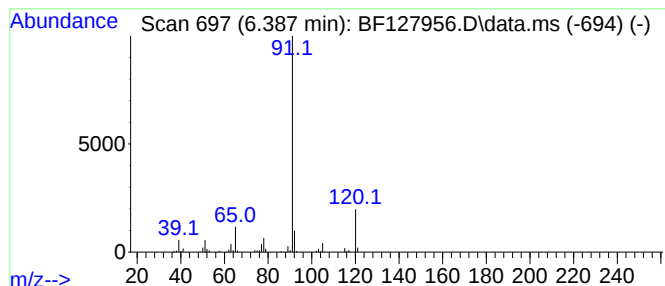
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Benzene, propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.387	3.85 ng	114919	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	94
2			N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78
3			1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	72
4			N-Isobutyl(phenyl)methanesulfona...	227	C11H17NO2S	144615-94-1	59
5			Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	53



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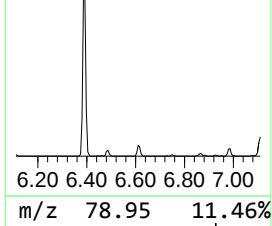
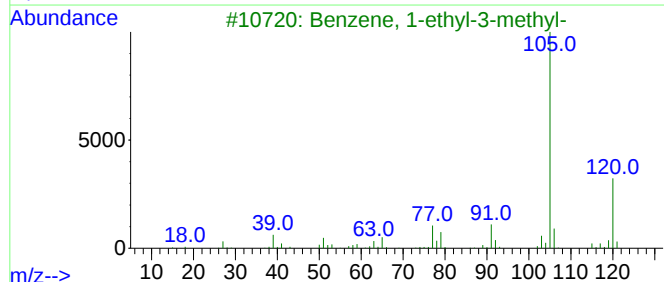
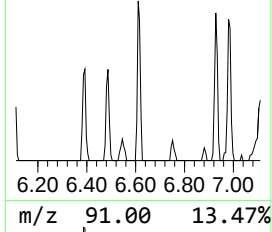
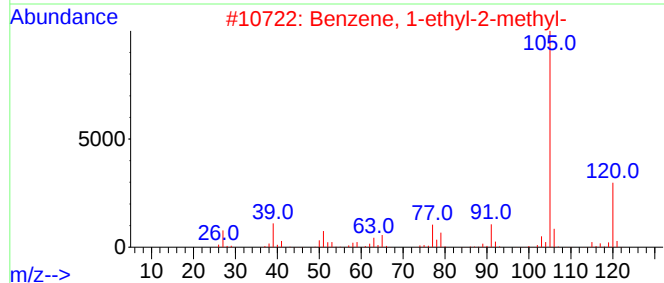
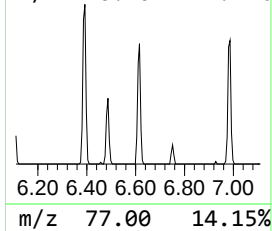
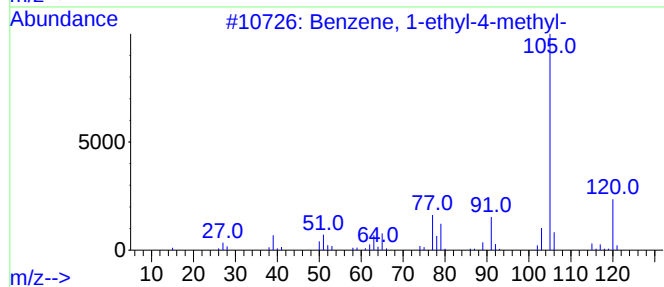
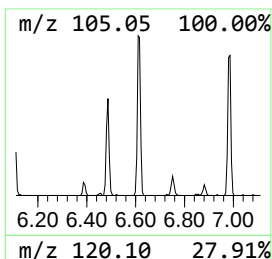
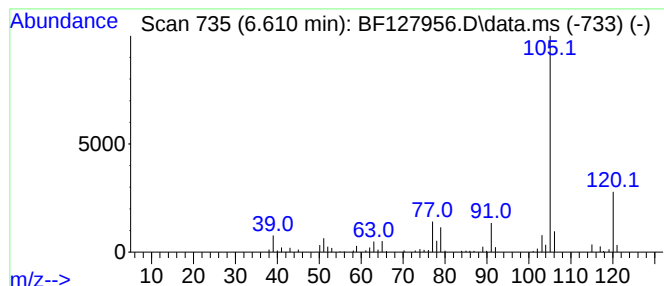
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-ethyl-4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.610	2.47 ng	73704	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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SPN-8

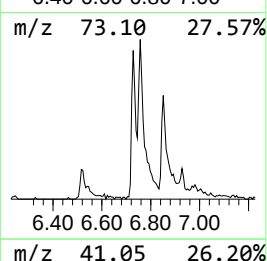
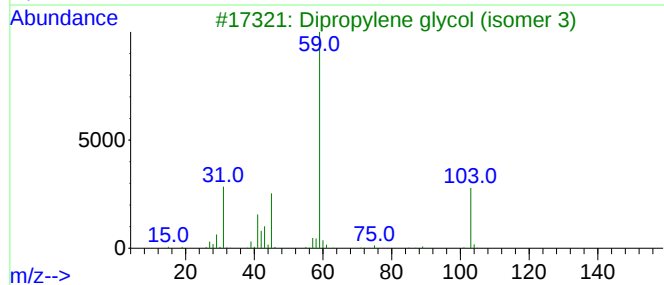
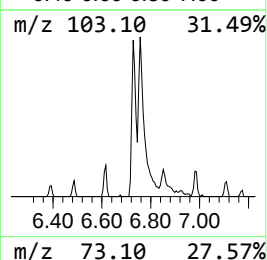
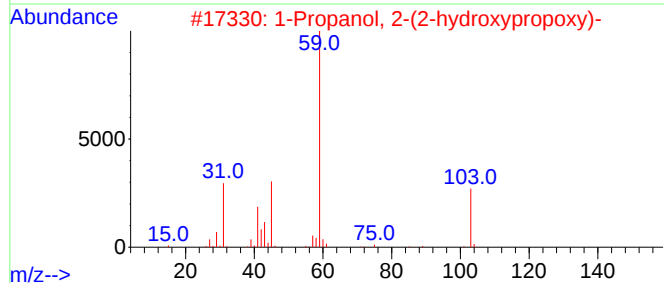
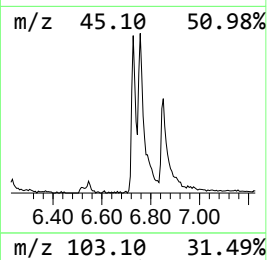
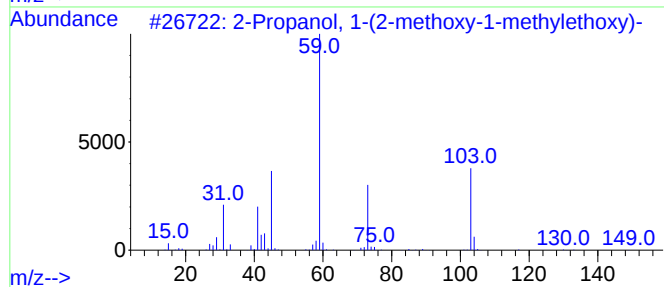
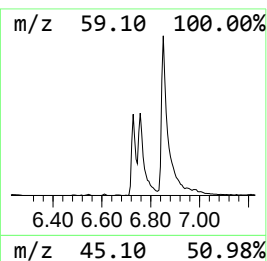
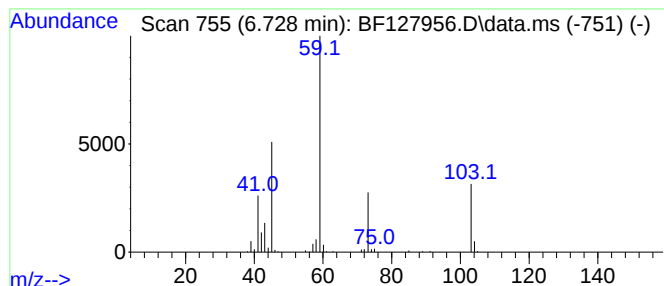
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.728	5.79 ng	172934	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		1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	59
3		Dipropylene glycol (isomer 3)	134	C6H14O3	1000506-25-5	53
4		1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	53
5		Dipropylene glycol (isomer 2)	134	C6H14O3	1000506-25-4	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042722\
Data File : BF127956.D
Acq On : 27 Apr 2022 20:24
Operator : CG\JU
Sample : N2482-15
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SPN-8

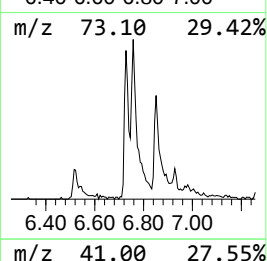
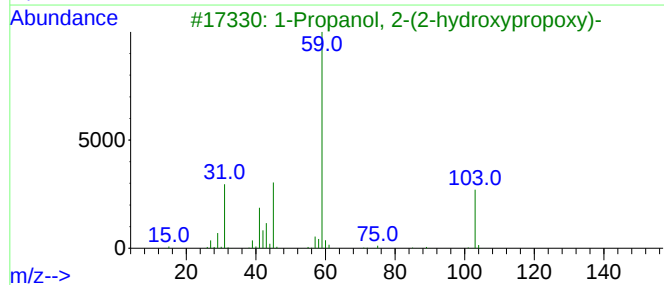
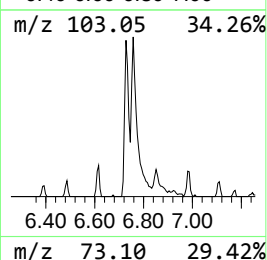
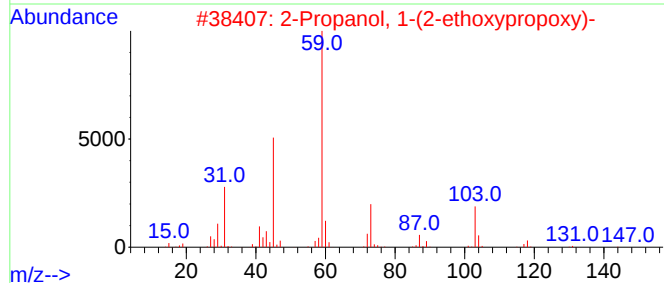
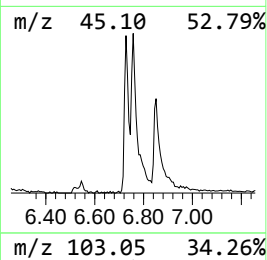
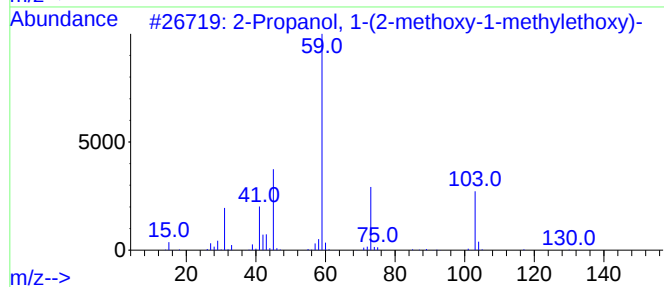
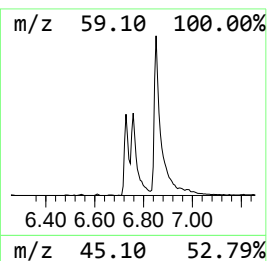
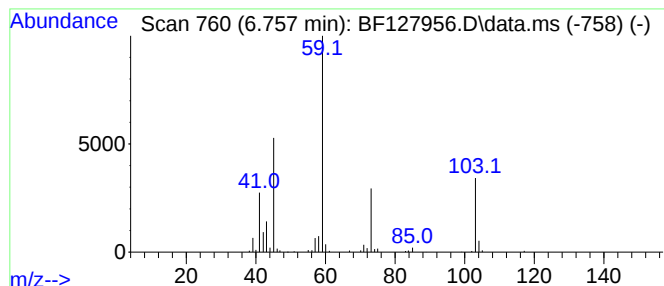
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 2-Propanol, 1-(2-ethoxyprop... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.757	6.77 ng	202186	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-ethoxypropoxy)-	162	C8H18O3	010143-32-5	72		
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	53		
4	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	50		
5	Dipropylene glycol (isomer 3)	134	C6H14O3	1000506-25-5	50		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042722\
Data File : BF127956.D
Acq On : 27 Apr 2022 20:24
Operator : CG\JU
Sample : N2482-15
Misc :
ALS Vial : 12 Sample Multiplier: 1

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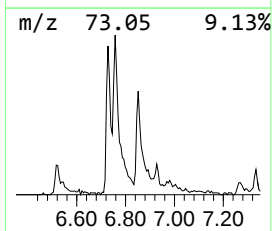
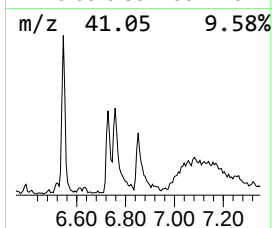
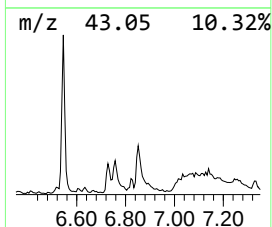
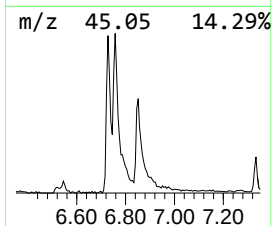
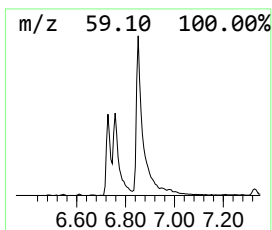
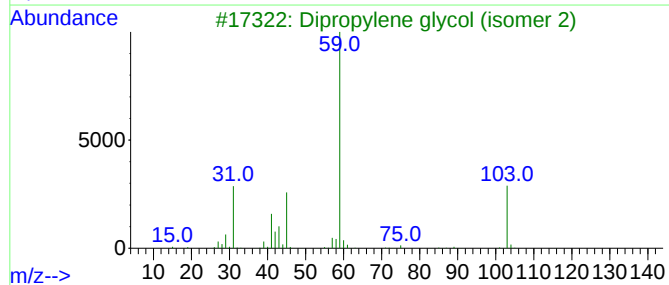
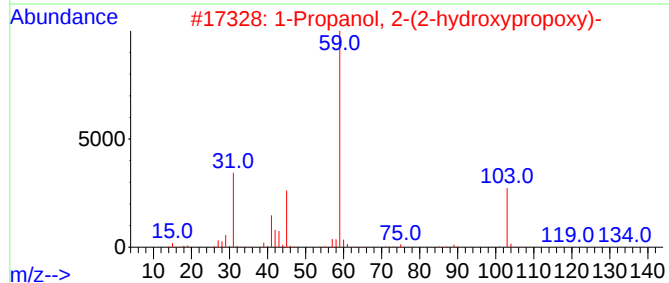
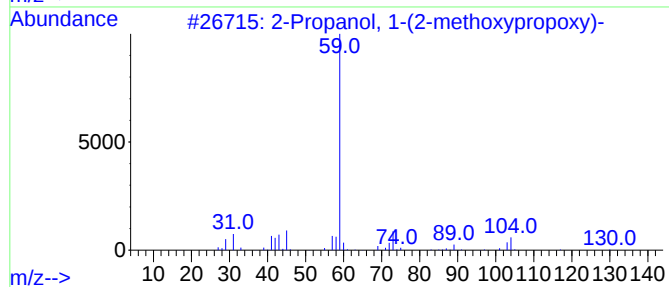
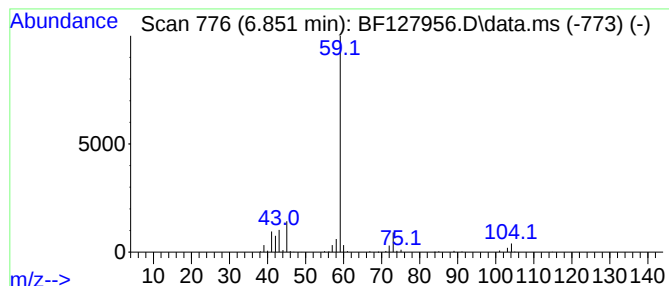
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 2-Propanol, 1-(2-methoxypro... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.851	8.89 ng	265614	1,4-Dichlorobenzene-d4	6.928

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	83
2		1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72
3		Dipropylene glycol (isomer 2)	134	C6H14O3	1000506-25-4	72
4		2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	64
5		Dipropylene glycol (isomer 3)	134	C6H14O3	1000506-25-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042722\
Data File : BF127956.D
Acq On : 27 Apr 2022 20:24
Operator : CG\JU
Sample : N2482-15
Misc :
ALS Vial : 12 Sample Multiplier: 1

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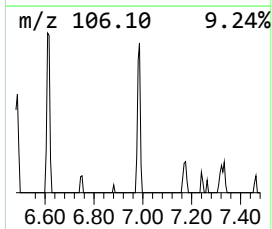
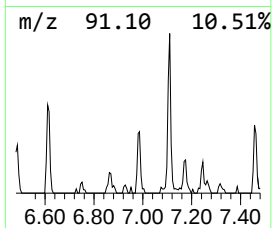
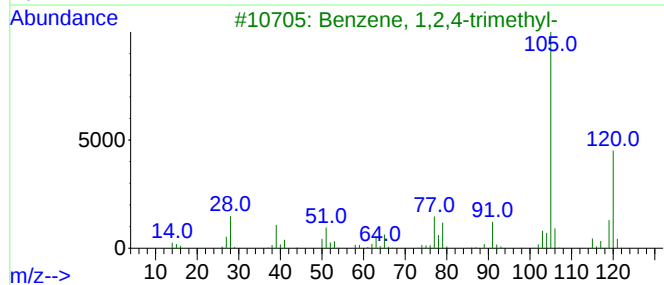
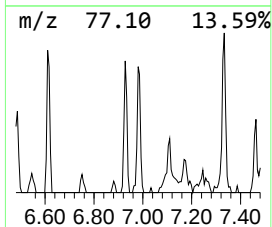
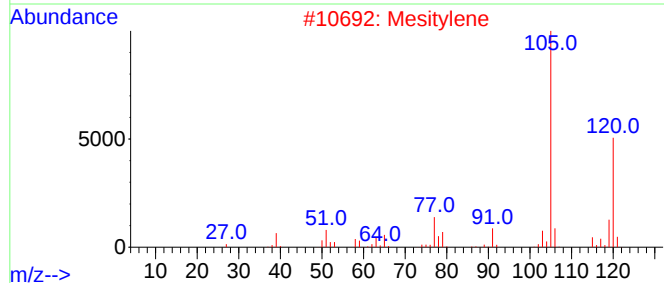
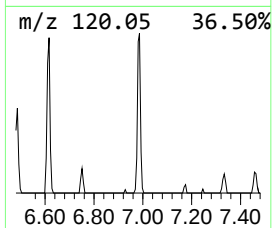
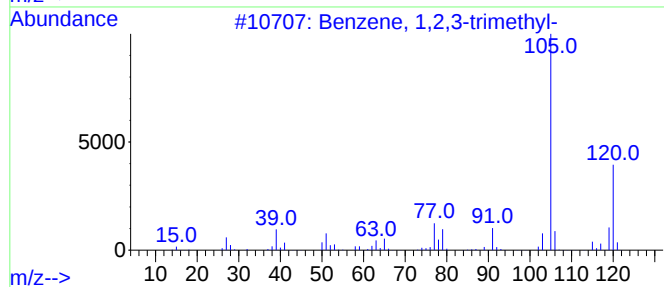
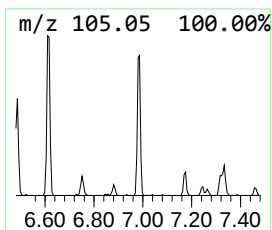
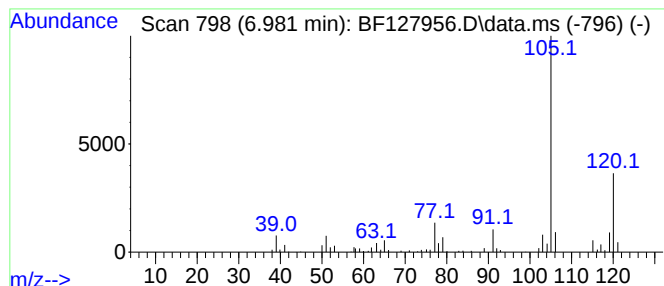
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.981	2.40 ng	71526	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2			Mesitylene	120	C9H12	000108-67-8	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042722\
Data File : BF127956.D
Acq On : 27 Apr 2022 20:24
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Sample : N2482-15
Misc :
ALS Vial : 12 Sample Multiplier: 1

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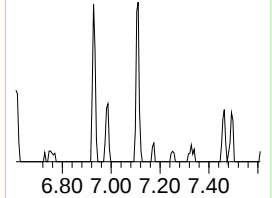
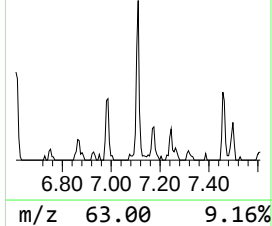
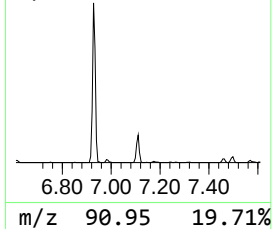
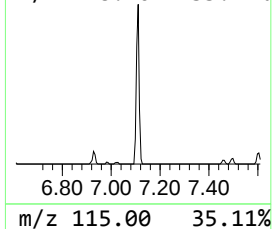
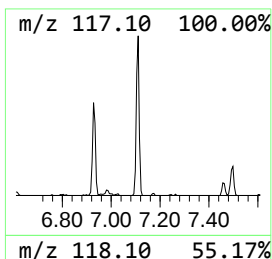
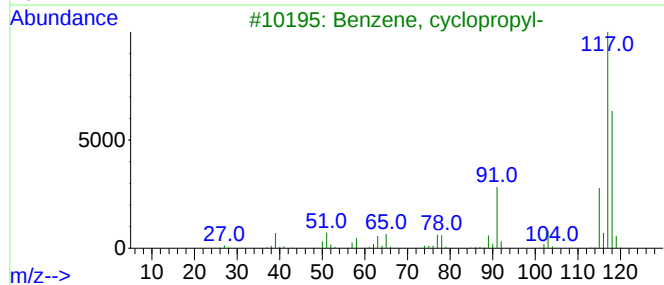
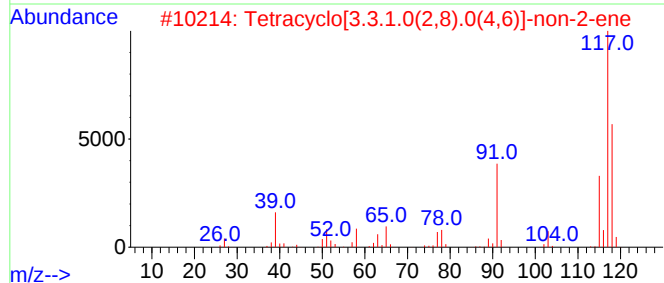
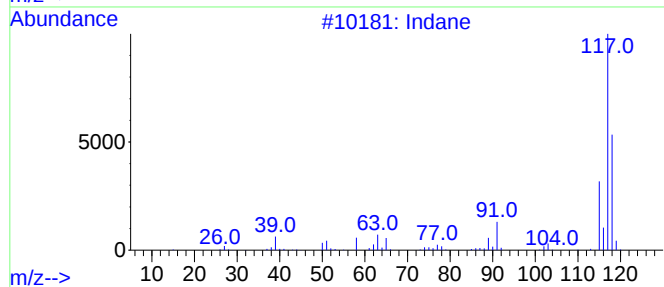
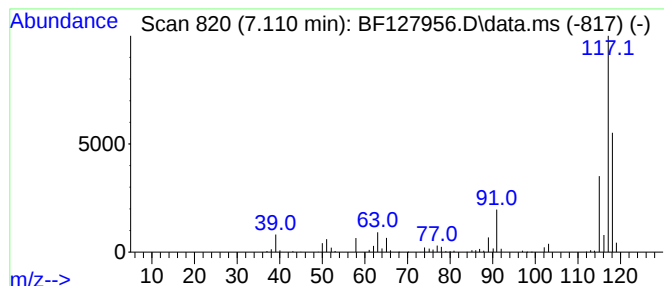
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Indane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.110	3.39 ng	101334	1,4-Dichlorobenzene-d4	6.928

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	94
2			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	72
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60
5			Benzene, 1-propenyl-	118	C9H10	000637-50-3	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042722\
Data File : BF127956.D
Acq On : 27 Apr 2022 20:24
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Misc :
ALS Vial : 12 Sample Multiplier: 1

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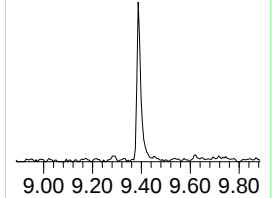
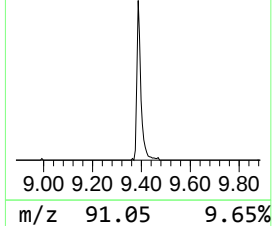
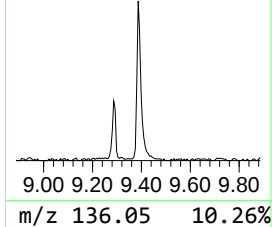
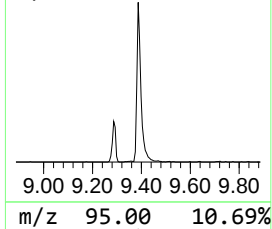
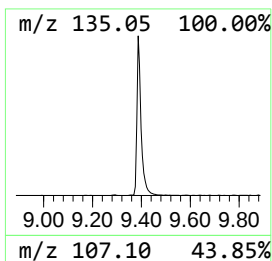
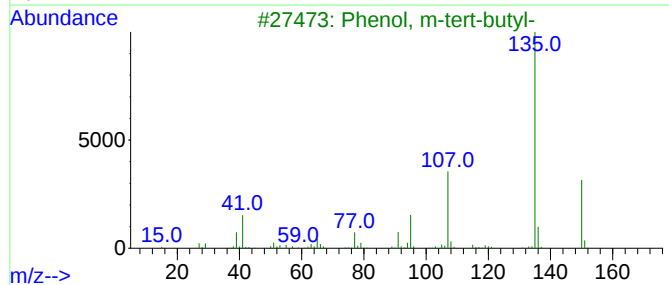
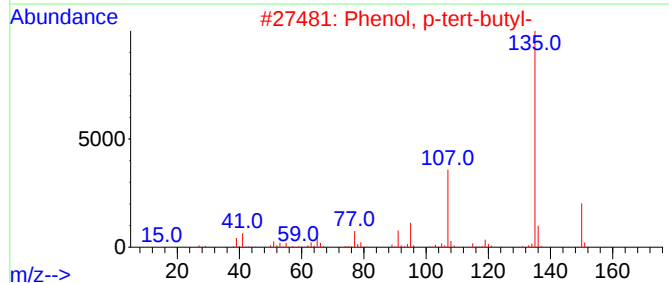
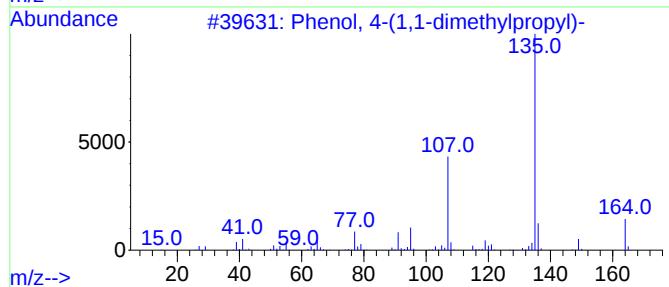
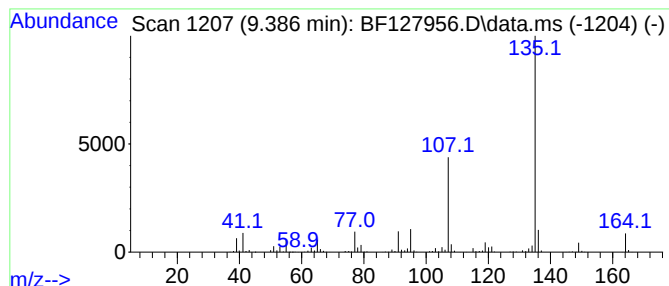
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Phenol, 4-(1,1-dimethylprop... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.386	14.34 ng	585356	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	90
3			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	64
4			4,4'-(1,2-diethylethylene)diphenol	270	C18H22O2	005635-50-7	64
5			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042722\
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Acq On : 27 Apr 2022 20:24
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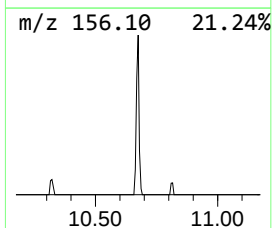
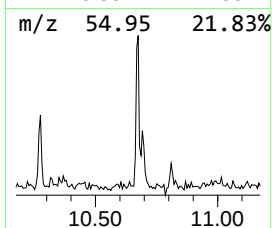
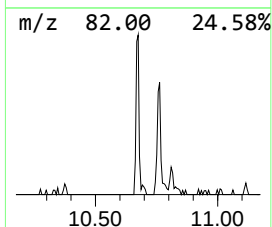
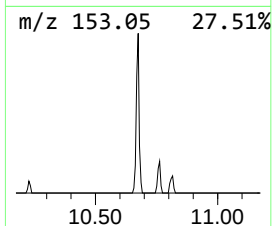
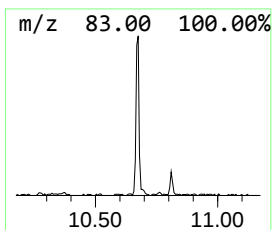
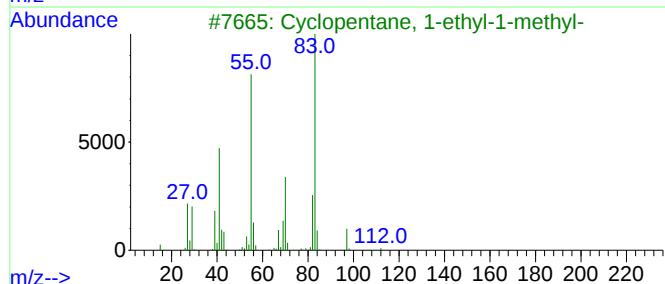
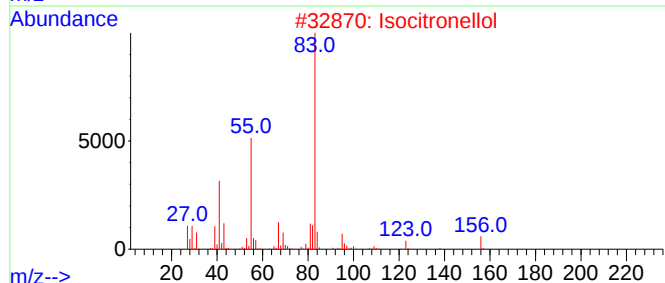
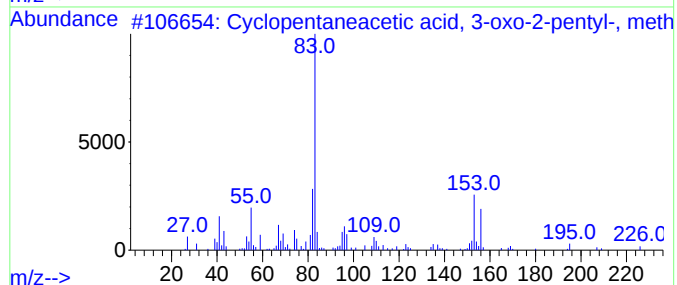
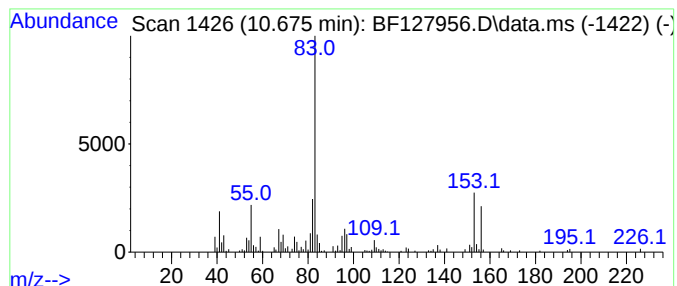
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Cyclopentaneacetic acid, 3-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.675	2.61 ng	106422	Acenaphthene-d10	9.969

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentaneacetic acid, 3-oxo-2...	226	C13H22O3	024851-98-7	95
2		Isocitronellol	156	C10H20O	018479-52-2	50
3		Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	47
4		N-(3-Piperidinyl)propanamide, N'...	226	C12H22N2O2	1000469-85-0	47
5		Oxalic acid, cyclohexyl tetradec...	368	C22H40O4	1000309-31-5	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042722\
Data File : BF127956.D
Acq On : 27 Apr 2022 20:24
Operator : CG\JU
Sample : N2482-15
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SPN-8

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
Cyclopentanone	4.457	13.5	ng	403347	1	6.928	597250	20.0
2-Propanol, 1-(...	5.157	4.8	ng	142487	1	6.928	597250	20.0
Benzene, (1-met...	6.104	2.2	ng	65013	1	6.928	597250	20.0
Benzene, propyl-	6.387	3.9	ng	114919	1	6.928	597250	20.0
Benzene, 1-ethy...	6.610	2.5	ng	73704	1	6.928	597250	20.0
2-Propanol, 1-(...	6.728	5.8	ng	172934	1	6.928	597250	20.0
2-Propanol, 1-(...	6.757	6.8	ng	202186	1	6.928	597250	20.0
2-Propanol, 1-(...	6.851	8.9	ng	265614	1	6.928	597250	20.0
Benzene, 1,2,3-...	6.981	2.4	ng	71526	1	6.928	597250	20.0
Indane	7.110	3.4	ng	101334	1	6.928	597250	20.0
Phenol, 4-(1,1-...	9.386	14.3	ng	585356	3	9.969	816431	20.0
Cyclopentaneace...	10.675	2.6	ng	106422	3	9.969	816431	20.0