

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081122\
 Data File : BF129726.D
 Acq On : 11 Aug 2022 21:07
 Operator : CG\JU
 Sample : N4024-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MANHOLE

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-BF072522.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF129726.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.393	185	188	203	rBV	143054	316333	5.96%	1.001%
2	5.469	538	541	556	rBV	1185021	1197382	22.55%	3.790%
3	6.134	650	654	662	rBV	154228	162837	3.07%	0.515%
4	6.481	710	713	721	rBV	633341	710704	13.39%	2.250%
5	6.828	769	772	776	rBV	1113781	953892	17.97%	3.019%
6	6.892	780	783	789	rBV	63485	77294	1.46%	0.245%
7	7.239	835	842	851	rBV2	241351	248682	4.68%	0.787%
8	7.398	858	869	872	rBV	2805516	3100515	58.40%	9.814%
9	7.657	907	913	918	rBV	44031	54753	1.03%	0.173%
10	8.022	971	975	982	rBV	233631	354037	6.67%	1.121%
11	8.110	986	990	994	rBV	1380844	1289211	24.28%	4.081%
12	8.275	1015	1018	1023	rBV	145382	244992	4.61%	0.775%
13	8.316	1023	1025	1029	rVB2	62007	69293	1.31%	0.219%
14	8.610	1071	1075	1077	rBV2	42922	53162	1.00%	0.168%
15	8.728	1090	1095	1096	rBV	46195	55424	1.04%	0.175%
16	8.745	1096	1098	1107	rVB2	53450	77354	1.46%	0.245%
17	9.063	1149	1152	1156	rBV2	46689	53284	1.00%	0.169%
18	9.192	1167	1174	1177	rBV	5252463	5309452	100.00%	16.806%
19	9.839	1281	1284	1286	rBV	139676	142640	2.69%	0.451%
20	9.869	1286	1289	1292	rVB	1805964	1423734	26.82%	4.507%
21	10.010	1308	1313	1315	rBV	116133	128998	2.43%	0.408%
22	10.051	1315	1320	1323	rVV	1466135	1891722	35.63%	5.988%
23	10.263	1352	1356	1358	rBV3	63806	69553	1.31%	0.220%
24	10.663	1419	1424	1427	rBV	2425490	2039462	38.41%	6.455%
25	10.786	1443	1445	1449	rVB	73893	67107	1.26%	0.212%
26	11.357	1537	1542	1546	rBV	1512328	1262286	23.77%	3.996%
27	11.539	1570	1573	1578	rBV	404231	326817	6.16%	1.034%
28	11.874	1627	1630	1634	rBV	128559	137672	2.59%	0.436%
29	12.663	1758	1764	1773	rBV	129468	166383	3.13%	0.527%
30	12.939	1807	1811	1815	rBV	2543344	2364379	44.53%	7.484%
31	13.868	1964	1969	1981	rVB5	83462	212600	4.00%	0.673%
32	14.004	1988	1992	1996	rBV	843670	745888	14.05%	2.361%
33	14.092	2002	2007	2012	rBV	117580	129922	2.45%	0.411%
34	14.615	2092	2096	2101	rBV	286775	262088	4.94%	0.830%
35	14.827	2129	2132	2135	rBV	56537	54485	1.03%	0.172%
36	15.474	2237	2242	2251	rVB	662083	871193	16.41%	2.758%

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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-BF072522.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	16.192	2356	2364	2385	rBV6	233504	1094175	20.61%	3.463%
38	16.357	2385	2392	2423	rVB3	923486	3453090	65.04%	10.930%
39	17.568	2588	2598	2620	rVB3	74651	419855	7.91%	1.329%

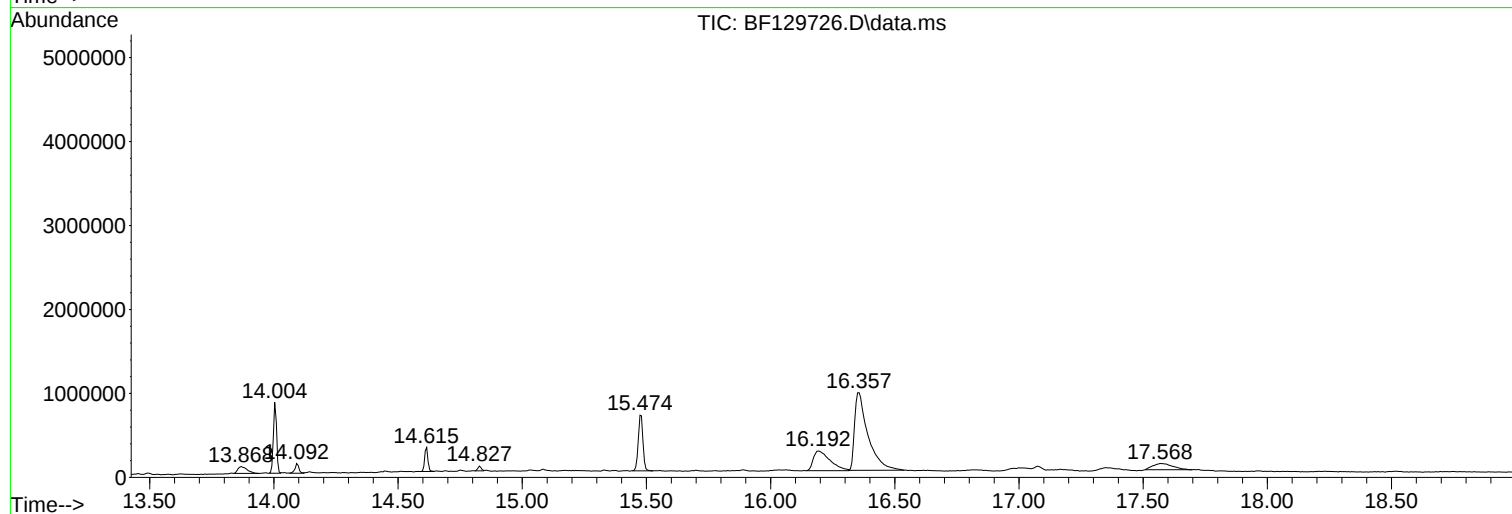
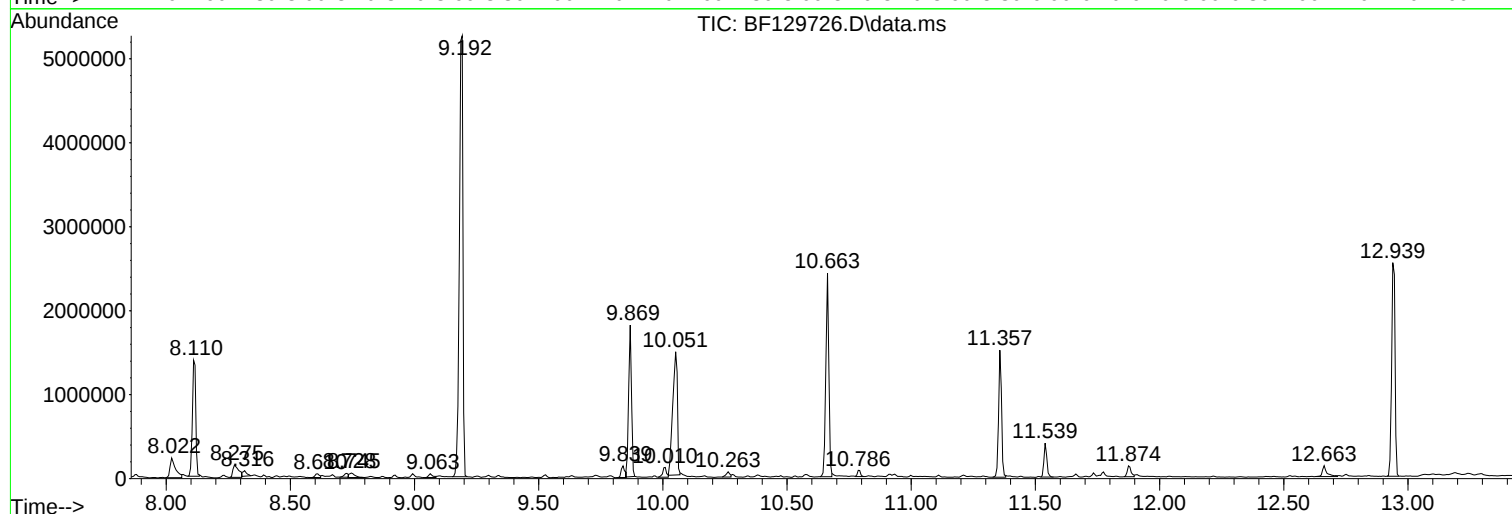
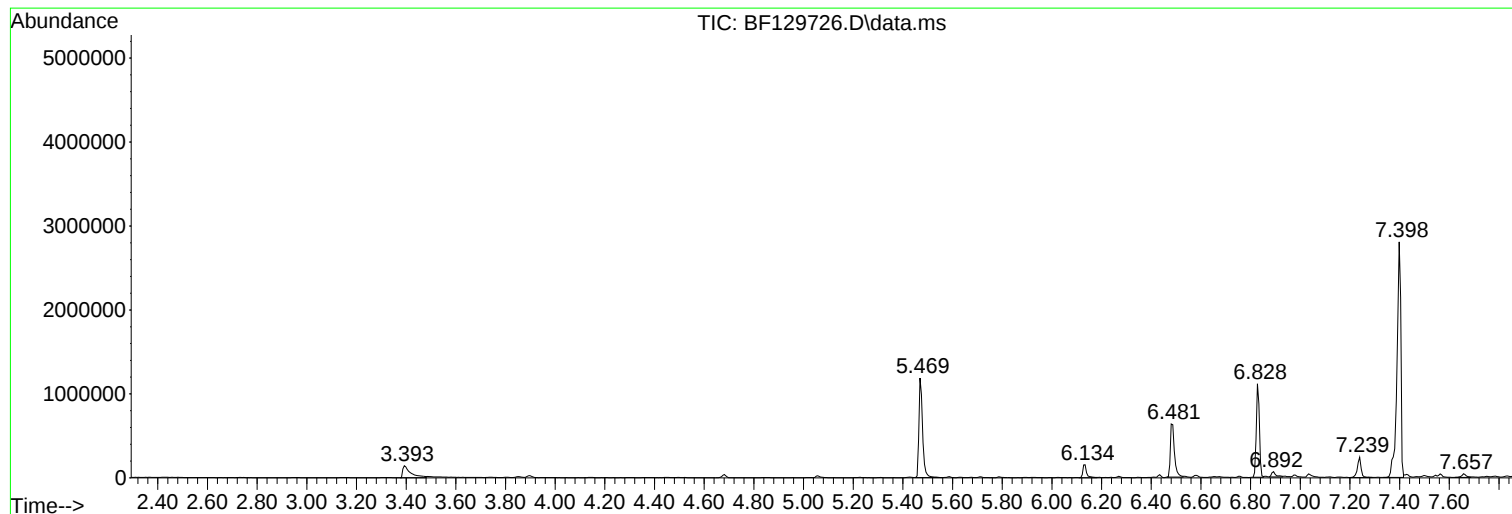
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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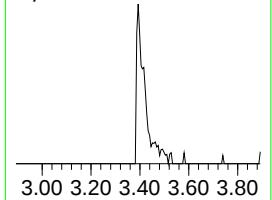
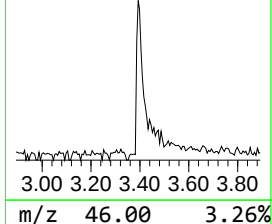
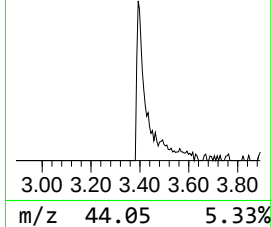
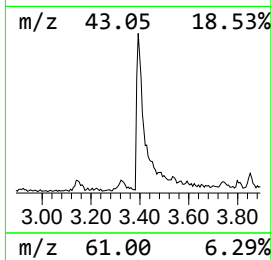
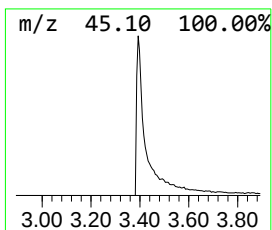
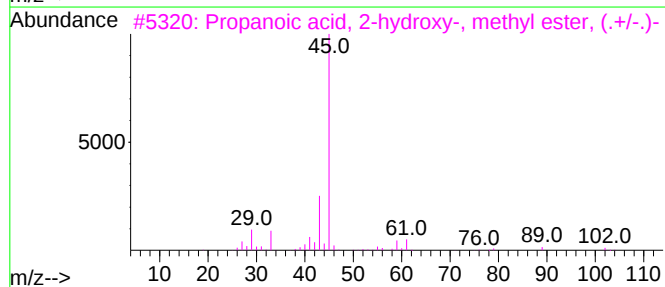
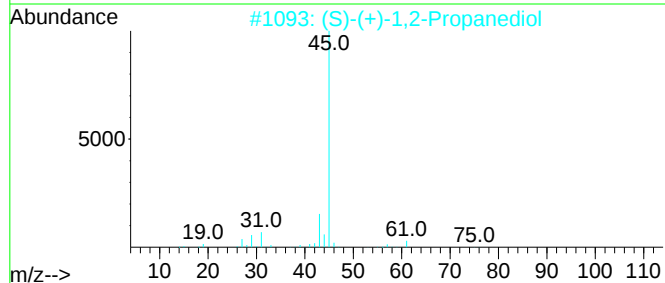
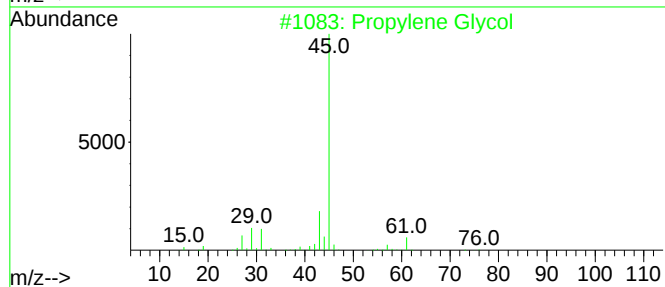
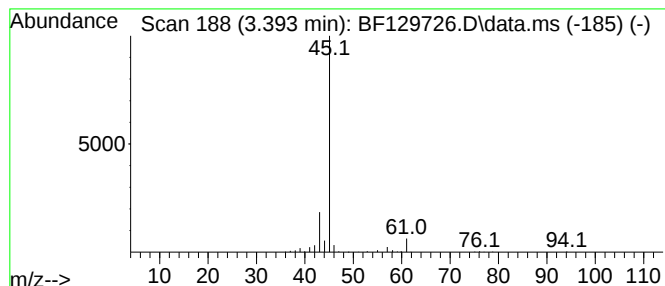
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.393	6.63 ng	316333	1,4-Dichlorobenzene-d4	6.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	90
2			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	56
3			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
4			Isopropyl Alcohol	60	C3H8O	000067-63-0	9
5			2-Hexanol	102	C6H14O	000626-93-7	9



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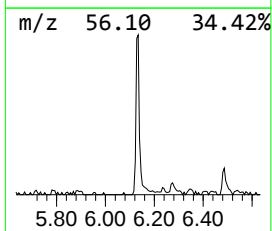
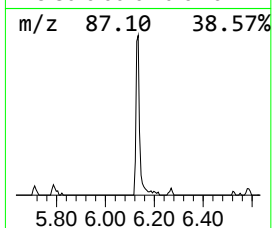
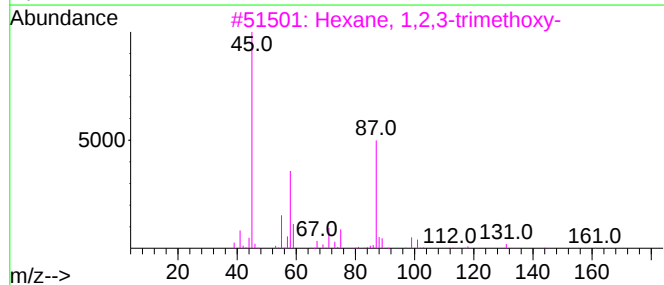
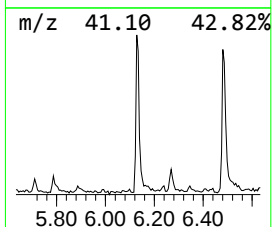
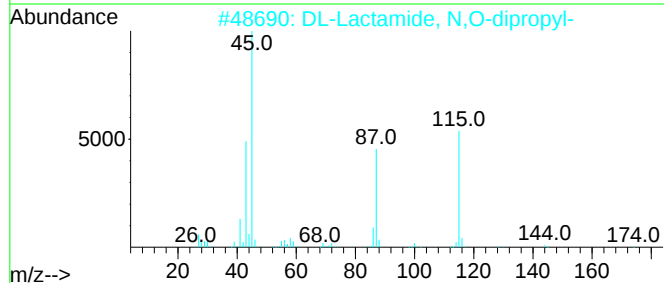
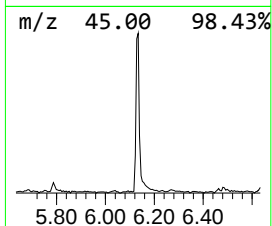
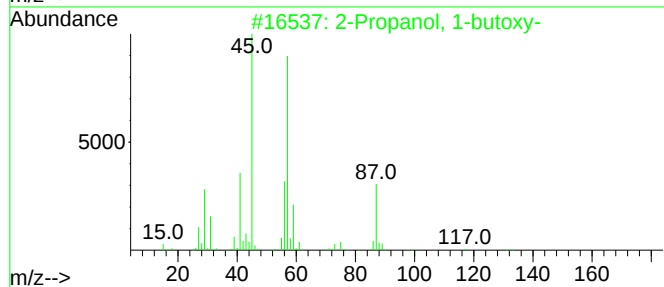
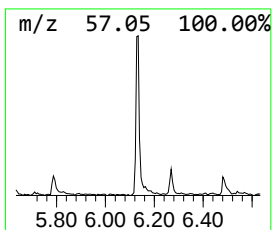
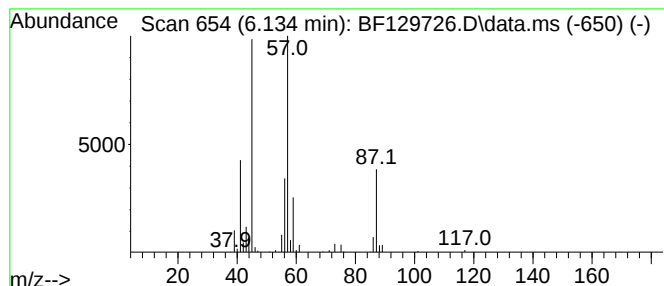
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Propanol, 1-butoxy- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.134	3.41 ng	162837	1,4-Dichlorobenzene-d4	6.828

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Propanol, 1-butoxy-	132	C7H16O2	005131-66-8	90
2		DL-Lactamide, N,O-dipropyl-	173	C9H19NO2	1010452-56-1	50
3		Hexane, 1,2,3-trimethoxy-	176	C9H20O3	020637-29-0	50
4		Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	47
5		1-Propanol, 2-(1-methylethoxy)-	118	C6H14O2	003944-37-4	45



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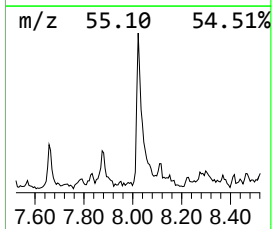
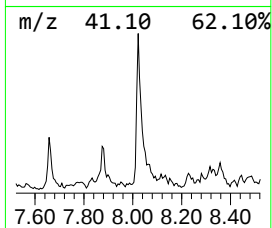
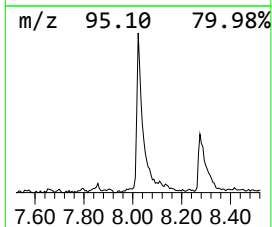
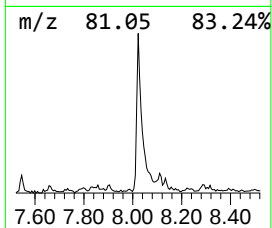
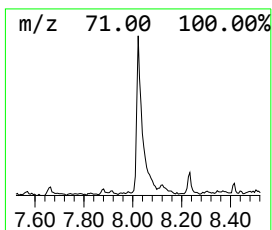
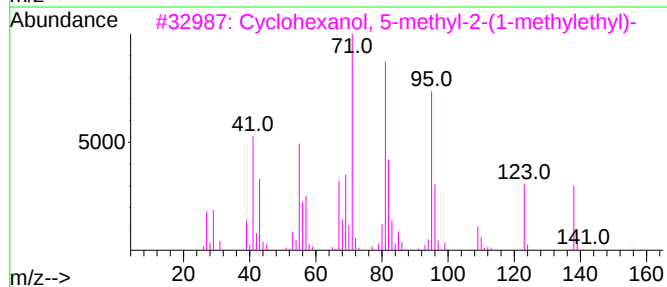
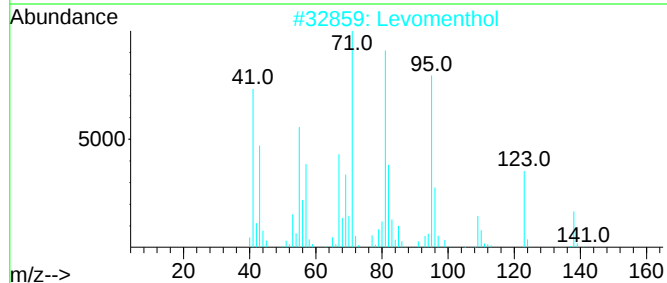
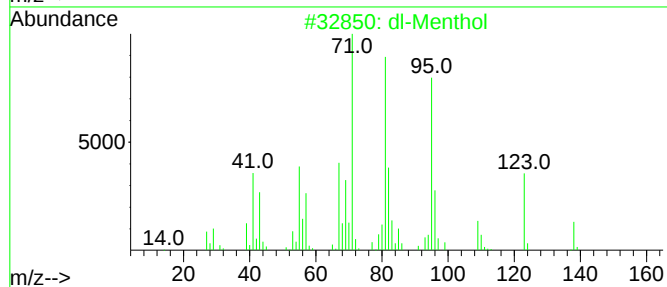
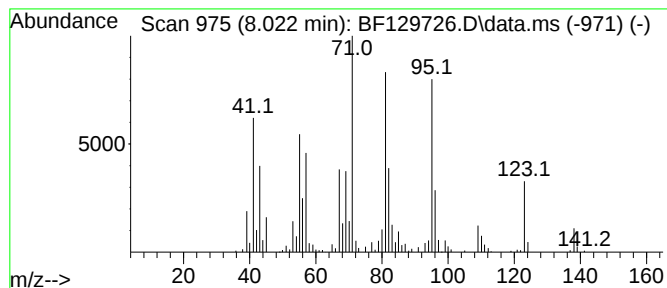
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 dl-Menthol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.022	5.49 ng	354037	Naphthalene-d8	8.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			dl-Menthol	156	C10H20O	000089-78-1	91
2			Levomenthol	156	C10H20O	002216-51-5	91
3			Cyclohexanol, 5-methyl-2-(1-meth...	156	C10H20O	001490-04-6	91
4			d-Menthol	156	C10H20O	015356-60-2	91
5			Cyclohexanol, 1-methyl-4-(1-meth...	156	C10H20O	021129-27-1	81



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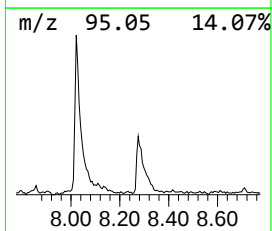
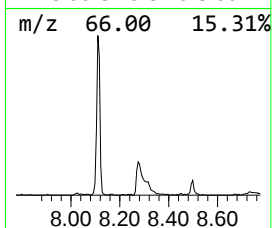
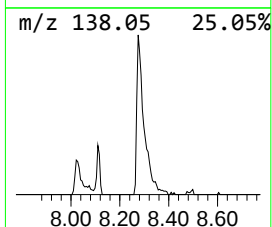
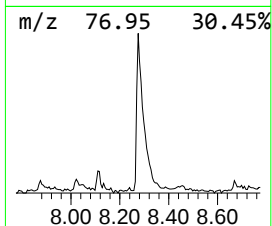
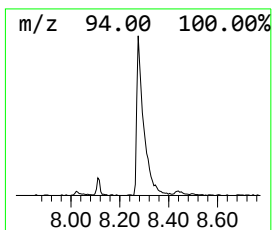
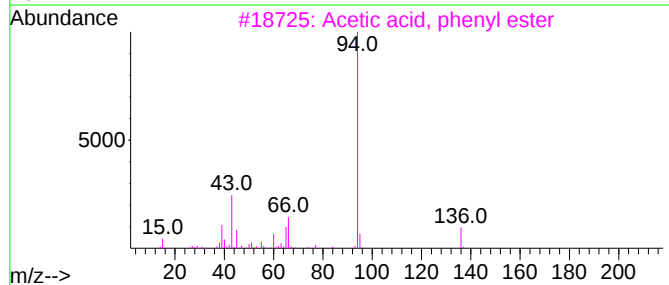
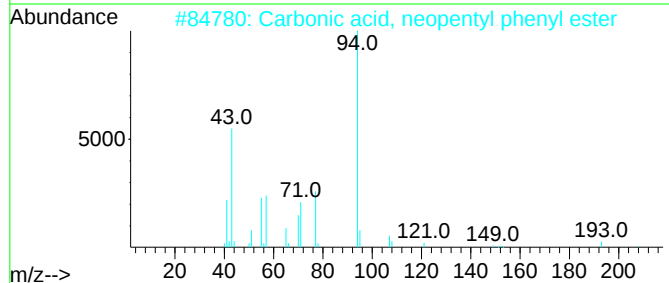
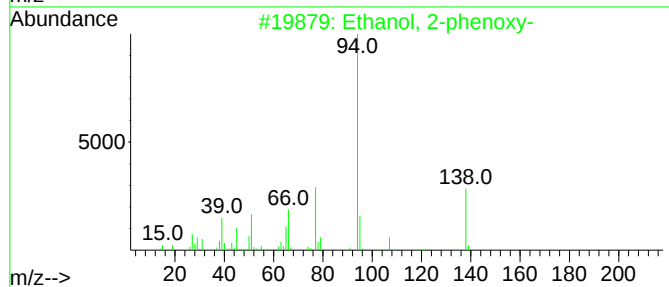
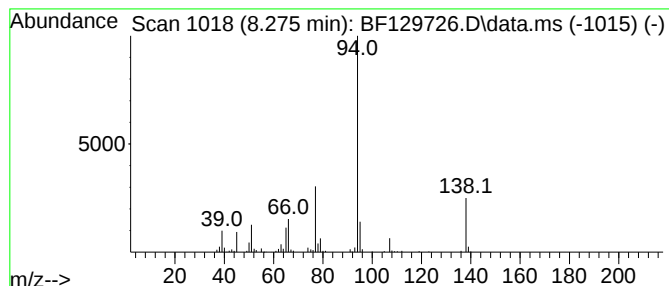
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Ethanol, 2-phenoxy- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.275	3.80 ng	244992	Naphthalene-d8	8.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	96
2			Carbonic acid, neopentyl phenyl ...	208	C12H16O3	013183-19-2	59
3			Acetic acid, phenyl ester	136	C8H8O2	000122-79-2	58
4			butanedioic acid, monophenyl ester	194	C10H10O4	1000400-29-4	53
5			Benzene, ethoxy-	122	C8H10O	000103-73-1	53



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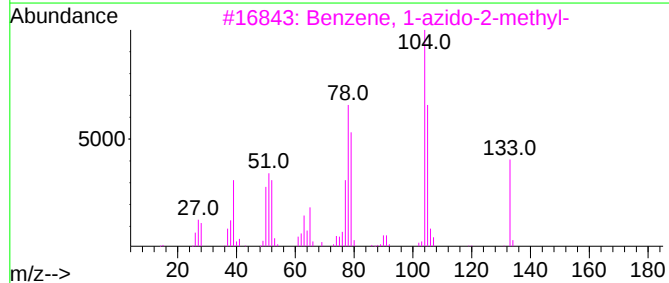
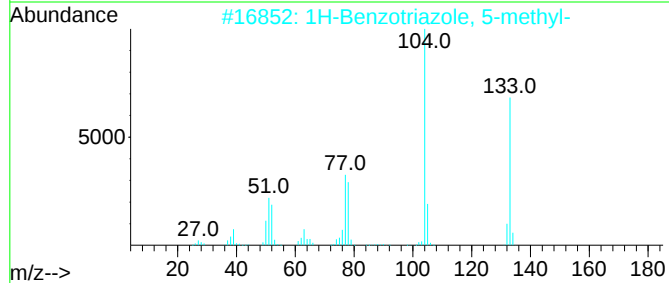
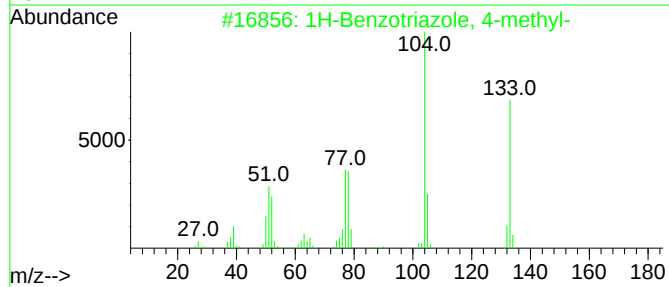
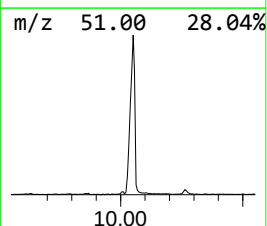
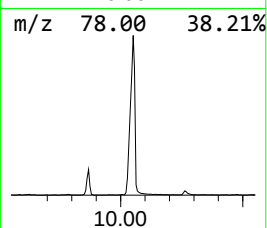
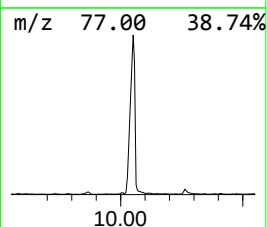
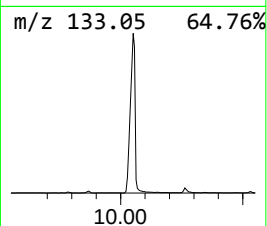
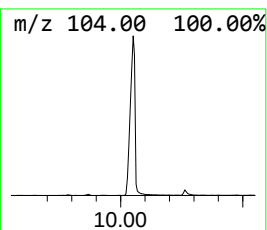
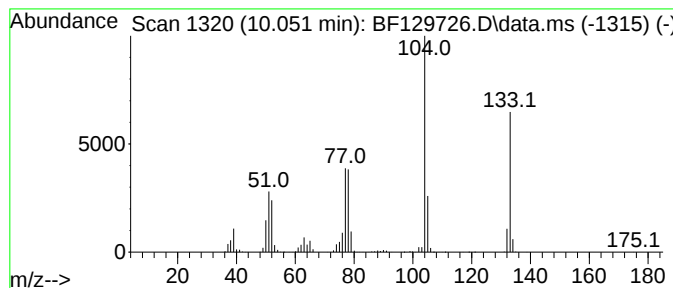
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 1H-Benzotriazole, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.051	26.57 ng	1891720	Acenaphthene-d10	9.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Benzotriazole, 4-methyl-	133	C7H7N3	029878-31-7	98
2			1H-Benzotriazole, 5-methyl-	133	C7H7N3	000136-85-6	97
3			Benzene, 1-azido-2-methyl-	133	C7H7N3	031656-92-5	72
4			2H-Indol-2-one, 1,3-dihydro-	133	C8H7NO	000059-48-3	72
5			Indan, epoxide	132	C9H8O	000768-22-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081122\
Data File : BF129726.D
Acq On : 11 Aug 2022 21:07
Operator : CG\JU
Sample : N4024-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MANHOLE

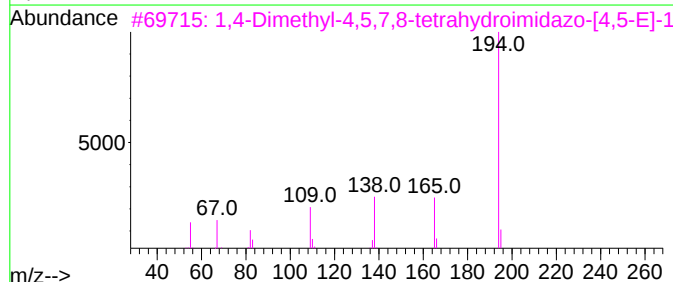
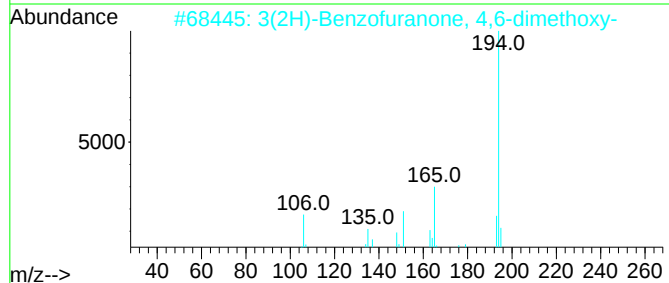
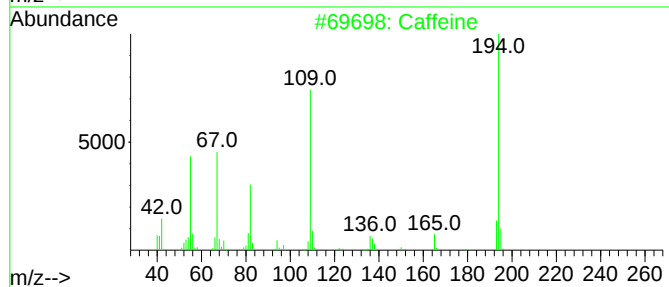
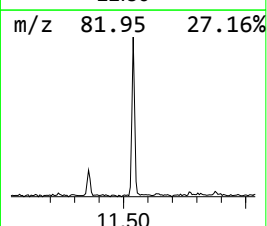
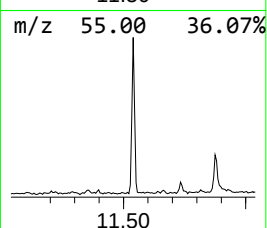
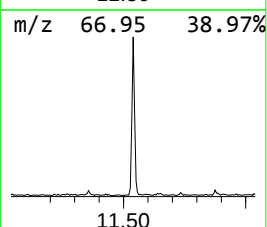
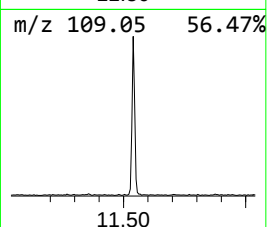
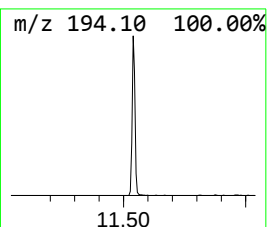
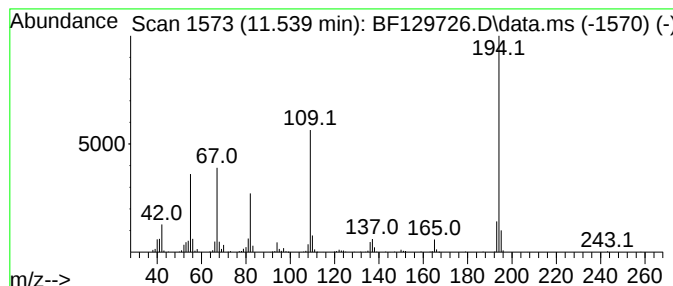
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Caffeine Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.539	5.18 ng	326817	Phenanthrene-d10	11.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Caffeine	194	C8H10N4O2	000058-08-2	98
2			3(2H)-Benzofuranone, 4,6-dimethoxy-	194	C10H10O4	004225-35-8	50
3			1,4-Dimethyl-4,5,7,8-tetrahydroi...	194	C8H10N4O2	130063-15-9	50
4			2H-Pyrazol-3-ol, 5-(2,5-dimethyl...	194	C9H10N2O5	1000304-22-5	43
5			2-Hydroxy-5,8-naphthalenedicarbo...	194	C12H6N2O	1000326-75-1	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081122\
Data File : BF129726.D
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Sample : N4024-01
Misc :
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ClientSampleId :
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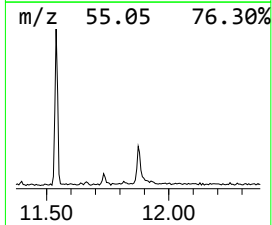
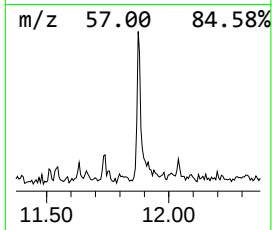
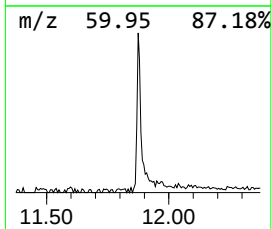
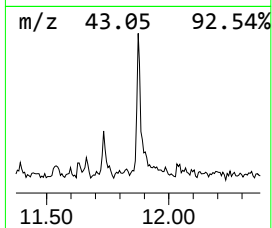
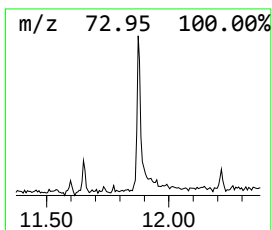
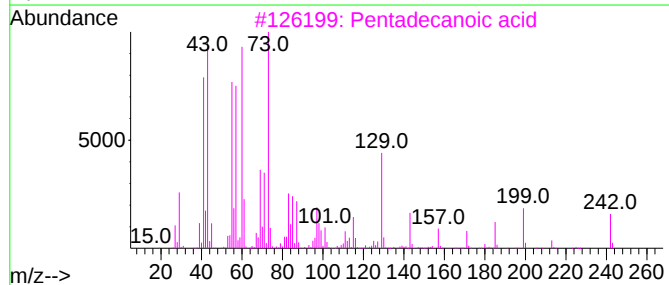
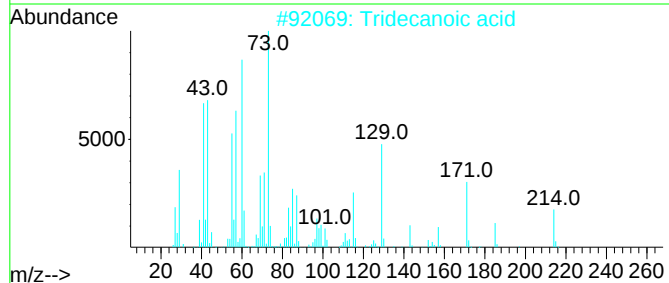
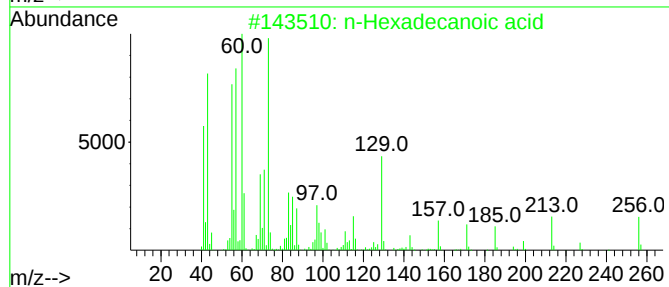
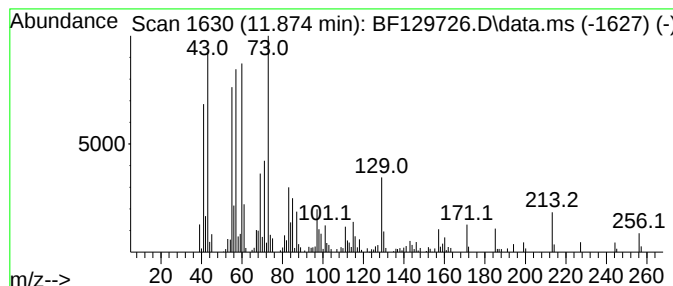
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.874	2.18 ng	137672	Phenanthrene-d10	11.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Tridecanoic acid	214	C13H26O2	000638-53-9	95
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5			Dodecanoic acid	200	C12H24O2	000143-07-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081122\
Data File : BF129726.D
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Misc :
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BNA_F
ClientSampleId :
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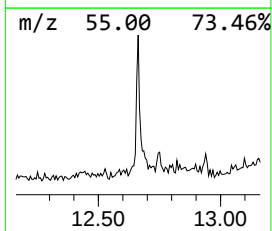
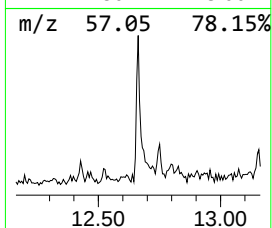
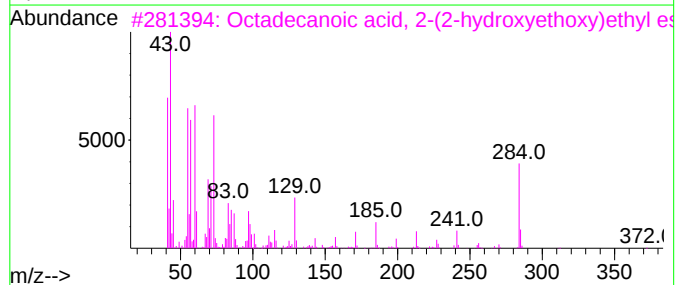
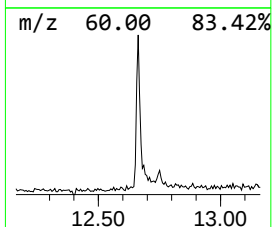
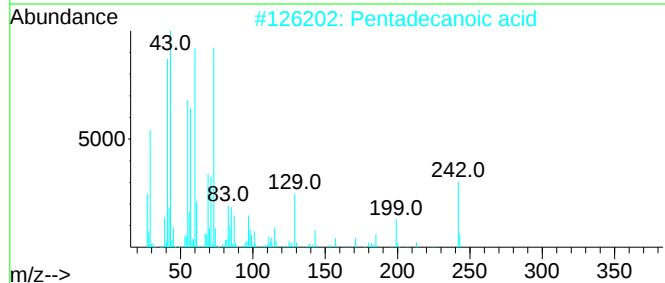
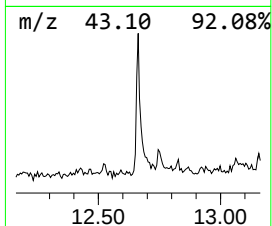
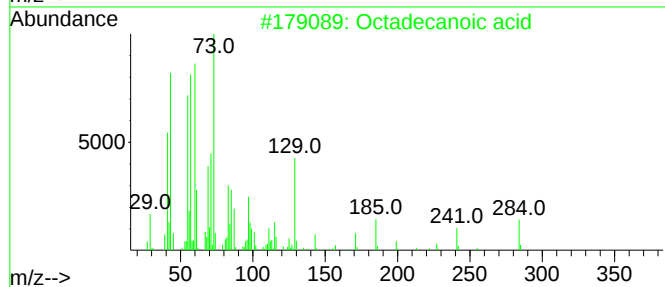
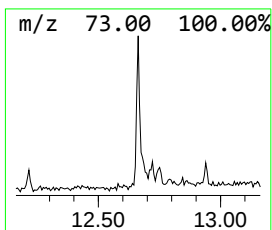
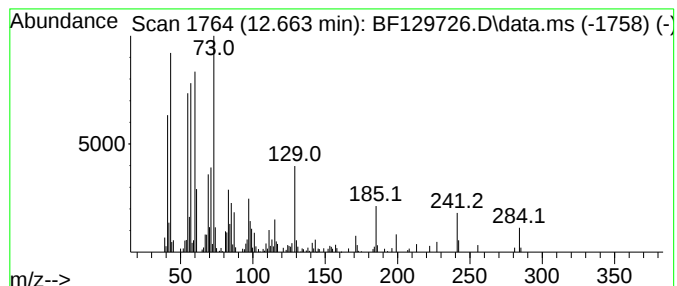
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.663	2.64 ng	166383	Phenanthrene-d10	11.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	87
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081122\
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Operator : CG\JU
Sample : N4024-01
Misc :
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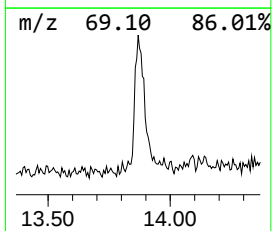
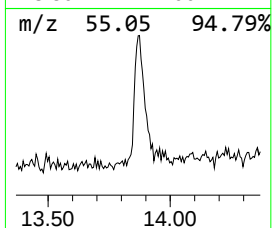
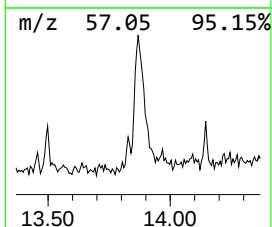
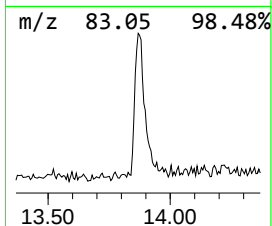
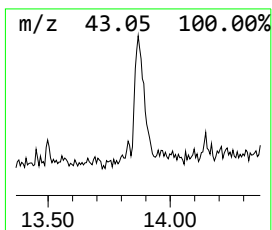
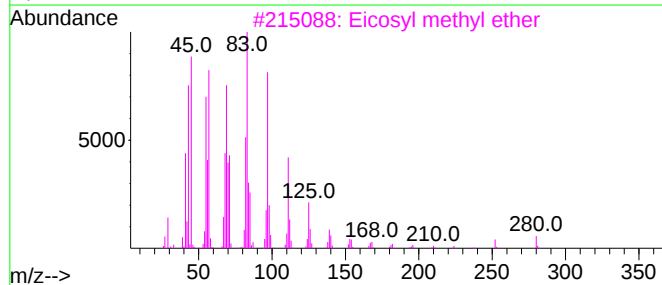
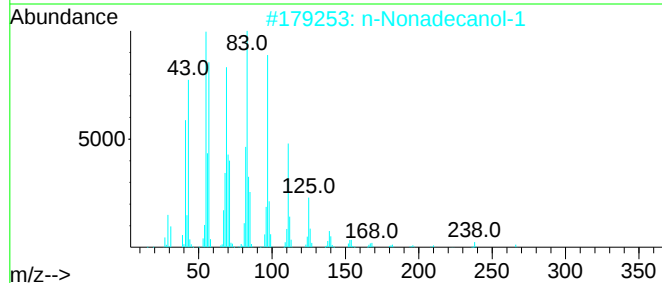
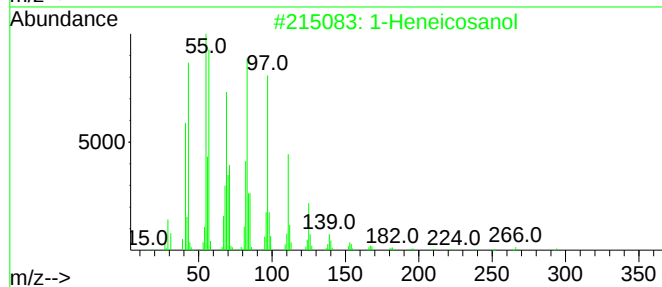
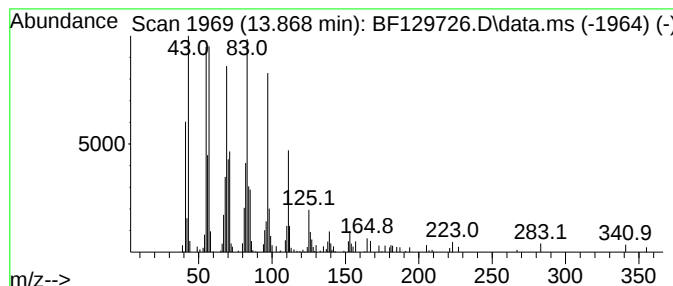
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Heneicosanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.868	5.70 ng	212600	Chrysene-d12	14.004	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	n-Nonadecanol-1	284	C19H40O	001454-84-8	95
3	Eicosyl methyl ether	312	C21H44O	1000406-29-4	95
4	1-Octadecanol	270	C18H38O	000112-92-5	95
5	Behenic alcohol	326	C22H46O	000661-19-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081122\
Data File : BF129726.D
Acq On : 11 Aug 2022 21:07
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ClientSampleId :
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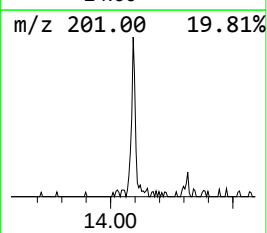
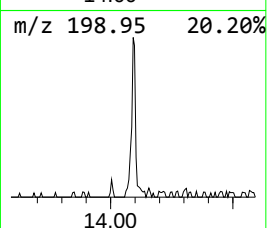
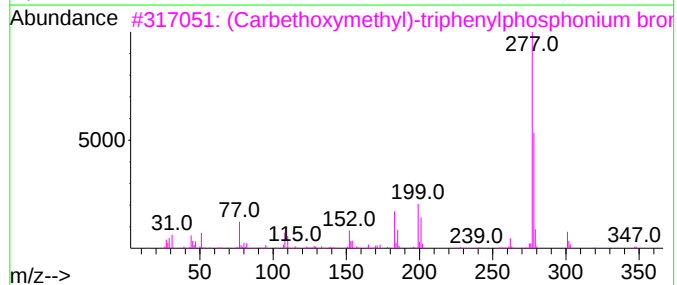
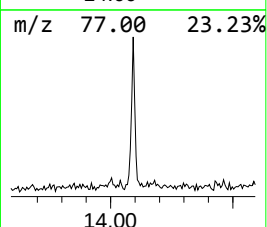
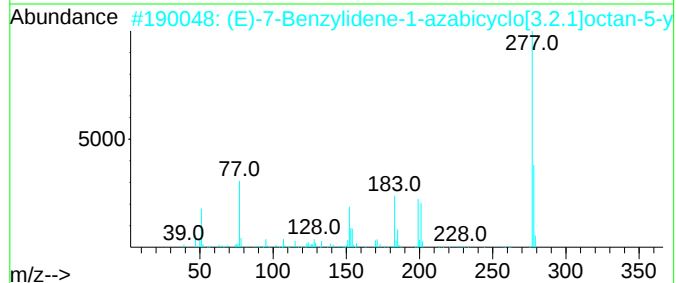
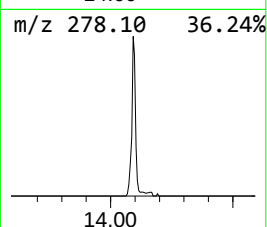
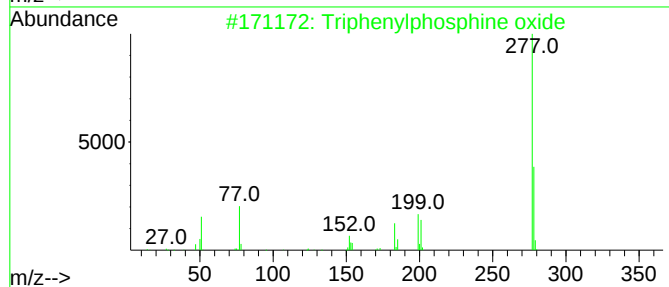
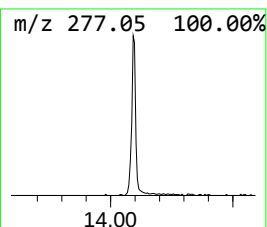
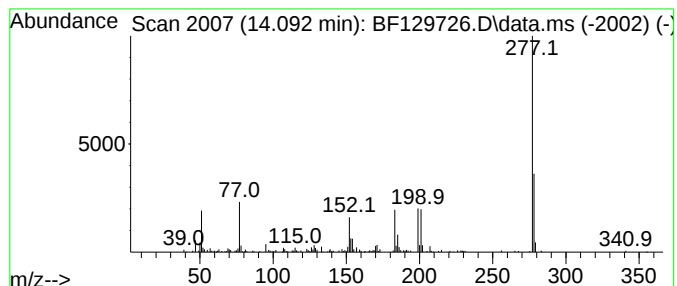
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Triphenylphosphine oxide Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.092	3.48 ng	129922	Chrysene-d12	14.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	97
2			(E)-7-Benzylidene-1-azabicyclo[3...]	293	C15H19NO3S	1000483-83-4	49
3			(Carbethoxymethyl)-triphenylphos...	428	C22H22BrO2P	001530-45-6	45
4			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	39
5			Cobalt,(.eta.<5>-2,4-cyclopentad...	277	C16H15BCo	036682-12-9	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081122\
Data File : BF129726.D
Acq On : 11 Aug 2022 21:07
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ALS Vial : 11 Sample Multiplier: 1

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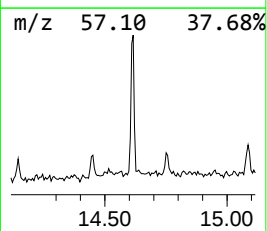
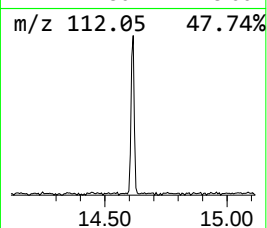
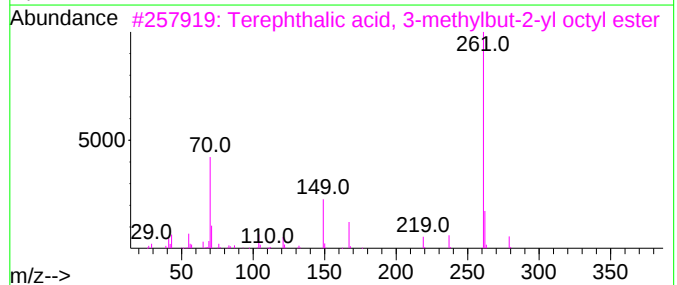
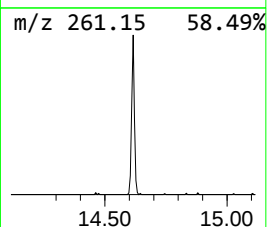
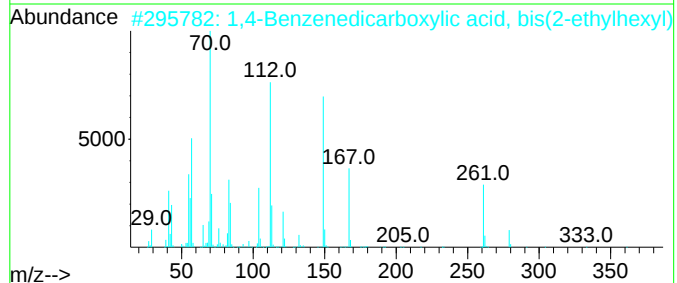
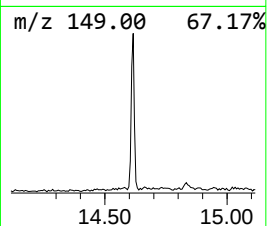
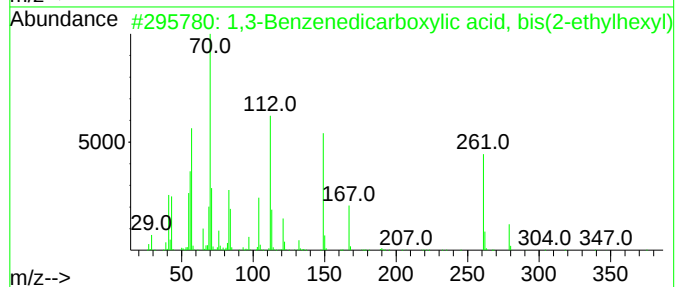
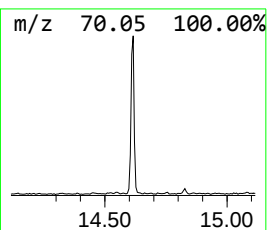
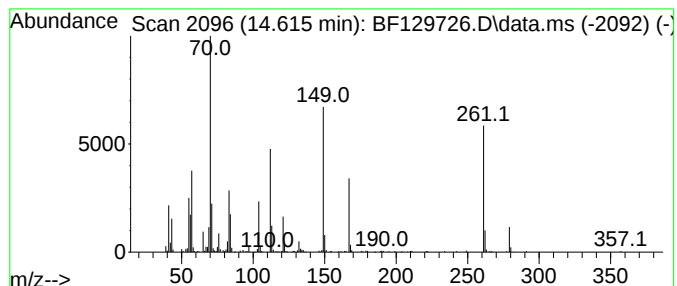
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1,3-Benzenedicarboxylic aci... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.616	7.03 ng	262088	Chrysene-d12	14.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	80
2			1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	70
3			Terephthalic acid, 3-methylbut-2...	348	C21H32O4	1000356-27-6	59
4			Terephthalic acid, 2-ethylhexyl ...	390	C24H38O4	1000324-00-5	53
5			Bromoacetic acid, 2-ethylhexyl e...	250	C10H19BrO2	068144-73-0	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081122\
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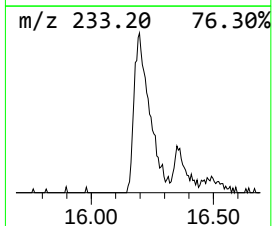
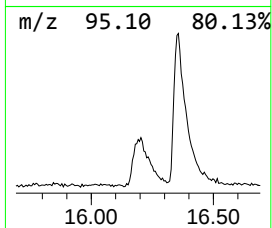
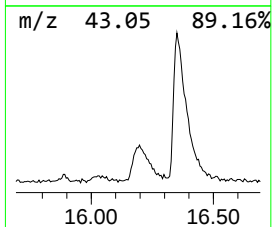
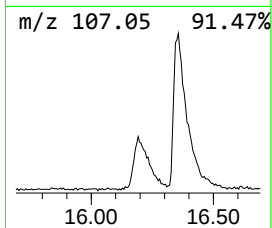
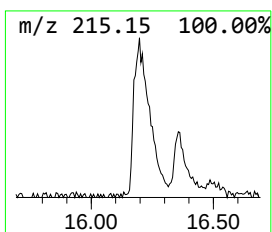
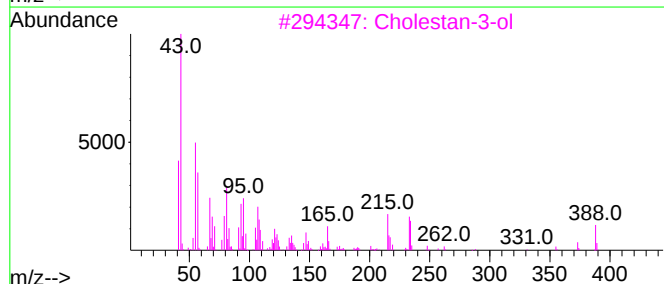
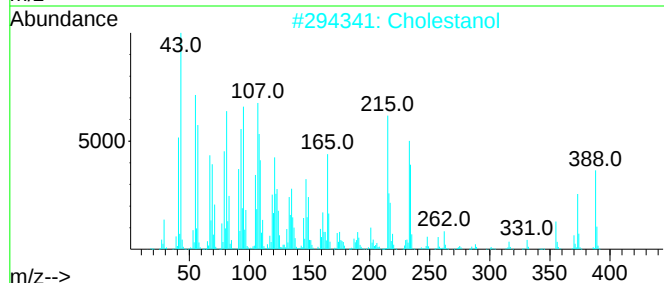
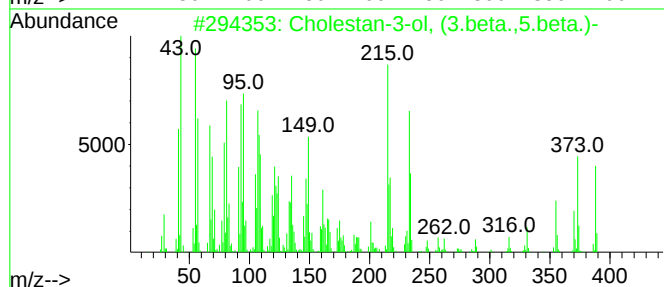
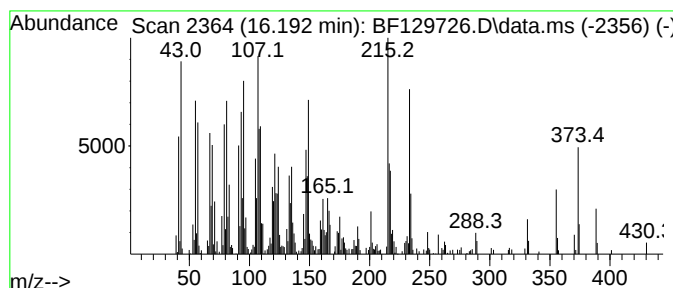
Instrument :
BNA_F
ClientSampleId :
MANHOLE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Cholestan-3-ol, (3.beta.,5.... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.192	25.12 ng	1094180	Perylene-d12	15.480	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cholestan-3-ol, (3.beta.,5.beta.)-	388	C27H48O	000360-68-9	90
2	Cholestanol	388	C27H48O	000080-97-7	81
3	Cholestan-3-ol	388	C27H48O	027409-41-2	70
4	trans-3-[2-Hydroxy-4-(2-methyl-2...	332	C22H36O2	1000370-41-7	38
5	Epicholestanol	388	C27H48O	000516-95-0	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081122\
Data File : BF129726.D
Acq On : 11 Aug 2022 21:07
Operator : CG\JU
Sample : N4024-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MANHOLE

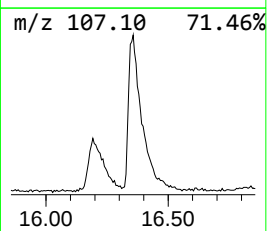
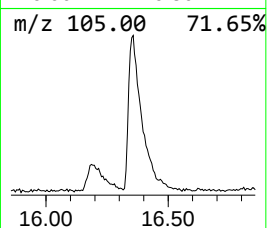
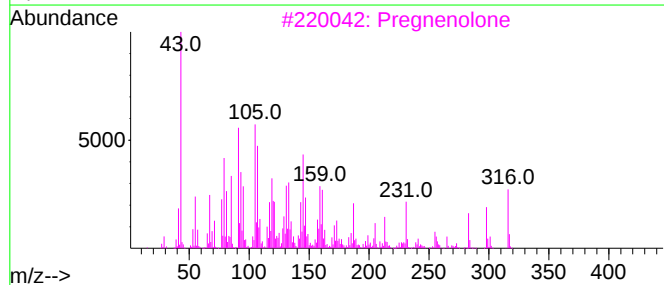
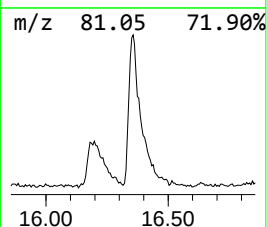
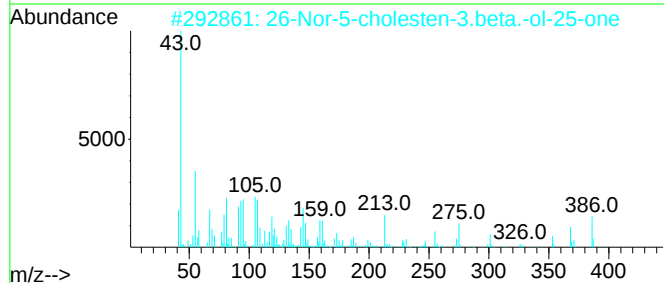
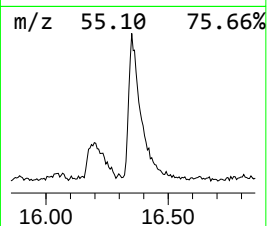
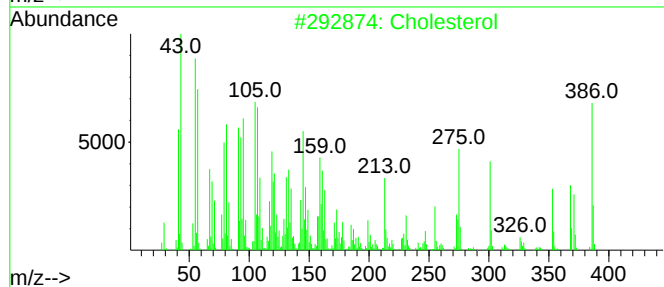
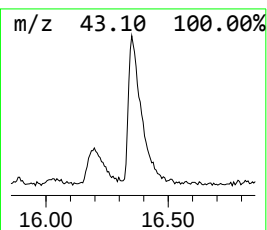
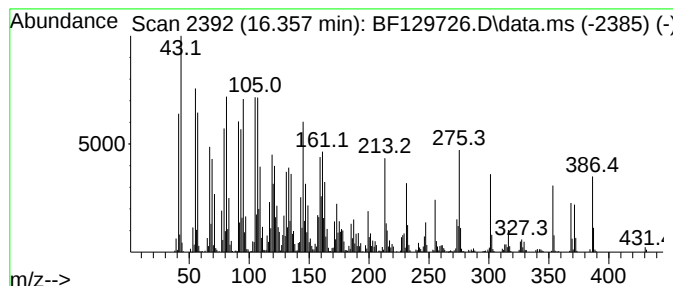
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Cholesterol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.357	79.27 ng	3453090	Perylene-d12	15.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cholesterol	386	C27H46O	000057-88-5	99
2			26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	96
3			Pregnenolone	316	C21H32O2	000145-13-1	56
4			Lathosterol	386	C27H46O	000080-99-9	50
5			Cholestane-3,5-diol, 5-acetate, ...	446	C29H50O3	041721-93-1	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081122\
Data File : BF129726.D
Acq On : 11 Aug 2022 21:07
Operator : CG\JU
Sample : N4024-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

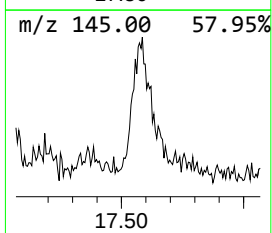
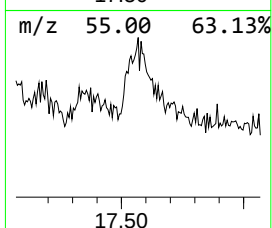
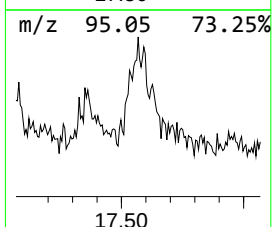
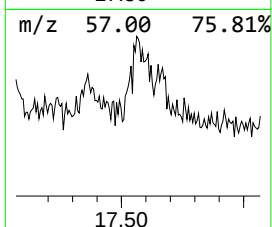
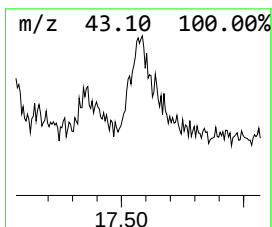
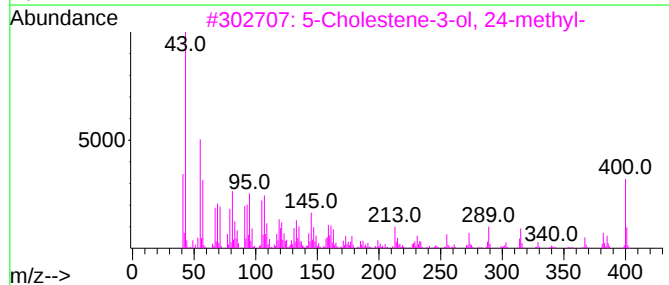
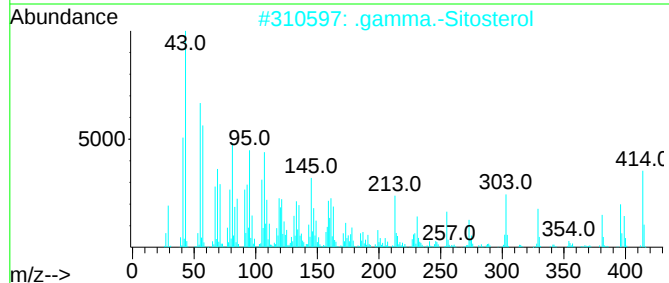
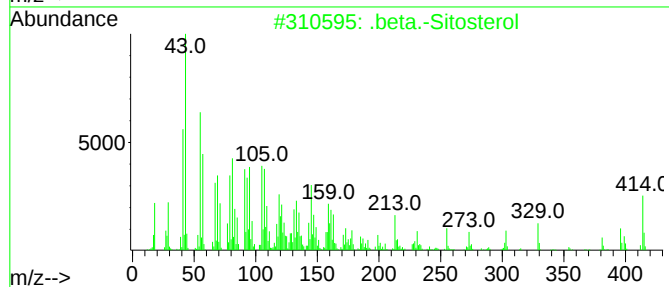
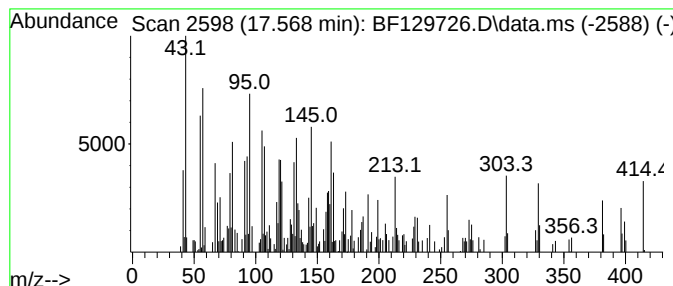
Instrument :
BNA_F
ClientSampleId :
MANHOLE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 .beta.-Sitosterol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.568	9.64 ng	419855	Perylene-d12		15.480	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	.beta.-Sitosterol		414	C29H50O	000083-46-5	78
2	.gamma.-Sitosterol		414	C29H50O	000083-47-6	64
3	5-Cholestene-3-ol, 24-methyl-		400	C28H48O	290299-12-6	42
4	Campesterol		400	C28H48O	000474-62-4	40
5	.beta.-Sitosterol acetate		456	C31H52O2	000915-05-9	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081122\
Data File : BF129726.D
Acq On : 11 Aug 2022 21:07
Operator : CG\JU
Sample : N4024-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MANHOLE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.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
Propylene Glycol	3.393	6.6	ng	316333	1	6.828	953892	20.0
2-Propanol, 1-b...	6.134	3.4	ng	162837	1	6.828	953892	20.0
dl-Menthol	8.022	5.5	ng	354037	2	8.110	1289210	20.0
Ethanol, 2-phen...	8.275	3.8	ng	244992	2	8.110	1289210	20.0
1H-Benzotriazol...	10.051	26.6	ng	1891720	3	9.869	1423730	20.0
Caffeine	11.539	5.2	ng	326817	4	11.357	1262290	20.0
n-Hexadecanoic ...	11.874	2.2	ng	137672	4	11.357	1262290	20.0
Octadecanoic acid	12.663	2.6	ng	166383	4	11.357	1262290	20.0
1-Heneicosanol	13.868	5.7	ng	212600	5	14.004	745888	20.0
Triphenylphosph...	14.092	3.5	ng	129922	5	14.004	745888	20.0
1,3-Benzenedica...	14.616	7.0	ng	262088	5	14.004	745888	20.0
Cholestan-3-ol,...	16.192	25.1	ng	1094180	6	15.480	871193	20.0
Cholesterol	16.357	79.3	ng	3453090	6	15.480	871193	20.0
.beta.-Sitosterol	17.568	9.6	ng	419855	6	15.480	871193	20.0