

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050624\
 Data File : BF137838.D
 Acq On : 06 May 2024 12:48
 Operator : CG\JU
 Sample : PB160665BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB160665BL

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-BF043024.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF137838.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.146	3	10	16	rVB	246326	304092	5.23%	0.970%
2	5.687	606	612	615	rBV	3537978	3684493	63.40%	11.758%
3	6.663	772	778	781	rBV	3139662	3651861	62.84%	11.654%
4	7.046	839	843	847	rVB	998842	836023	14.39%	2.668%
5	7.610	933	939	942	rBV	2294061	2424299	41.72%	7.736%
6	8.328	1056	1061	1065	rBV	1136925	1136877	19.56%	3.628%
7	9.410	1238	1245	1248	rBV	4872060	5042359	86.77%	16.091%
8	10.093	1355	1361	1364	rBV	1673035	1353181	23.29%	4.318%
9	10.887	1489	1496	1499	rBV	3515031	3197774	55.03%	10.205%
10	11.587	1611	1615	1618	rBV	1792816	1440722	24.79%	4.598%
11	13.181	1880	1886	1889	rBV	6054737	5811215	100.00%	18.545%
12	14.233	2061	2065	2069	rBV	1314311	1265744	21.78%	4.039%
13	15.804	2326	2332	2343	rBV	766193	987908	17.00%	3.153%
14	18.274	2736	2752	2765	rBV5	43177	199803	3.44%	0.638%

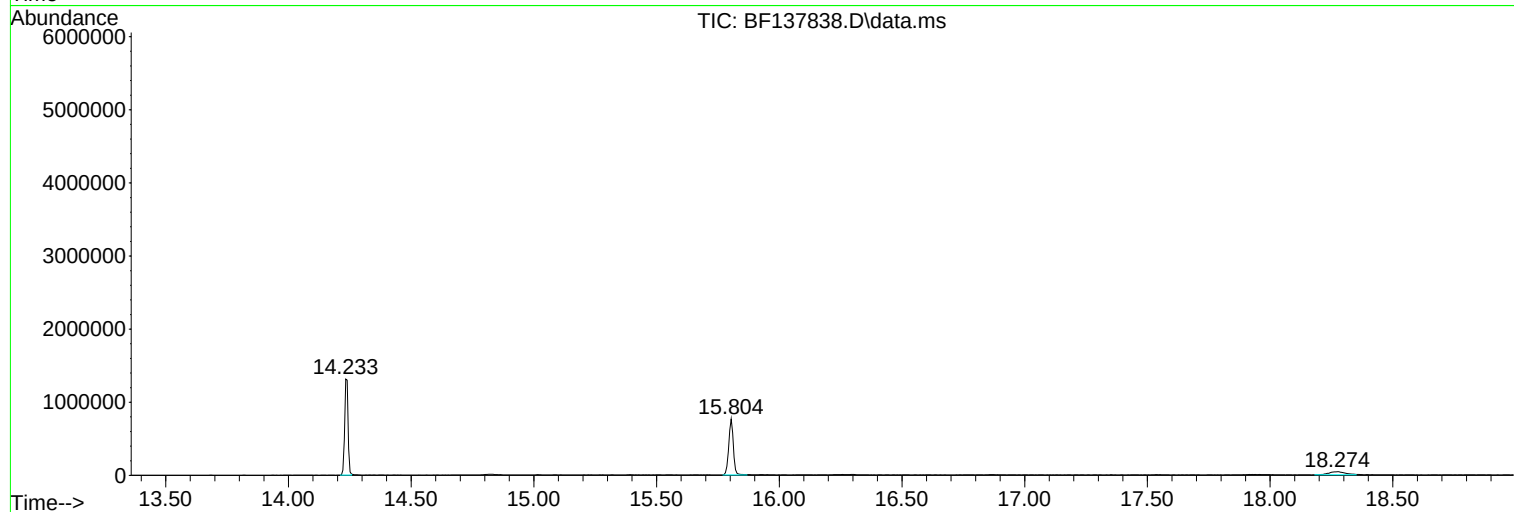
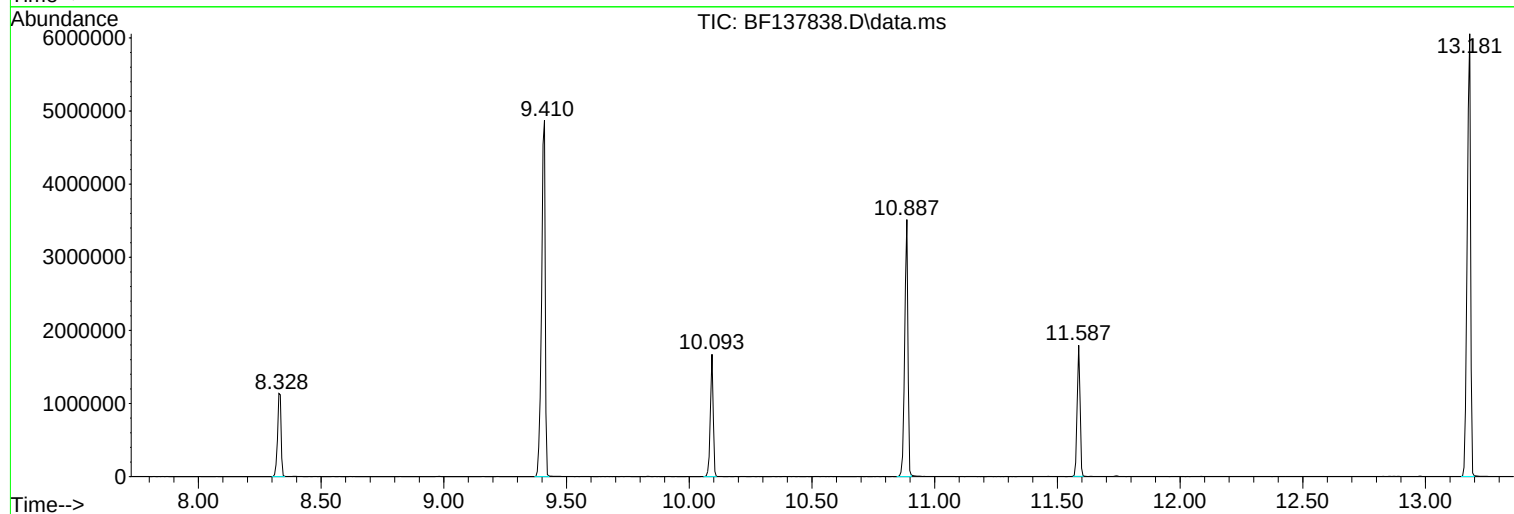
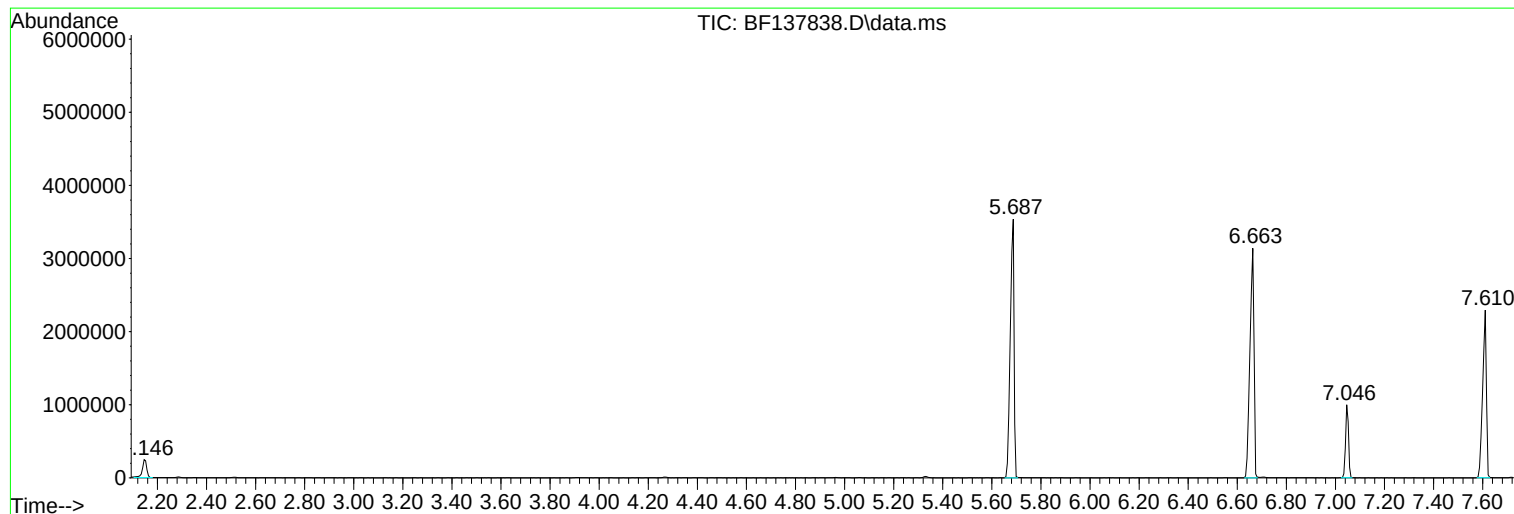
Sum of corrected areas: 31336351

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF043024.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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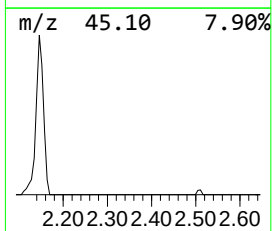
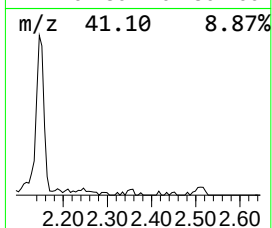
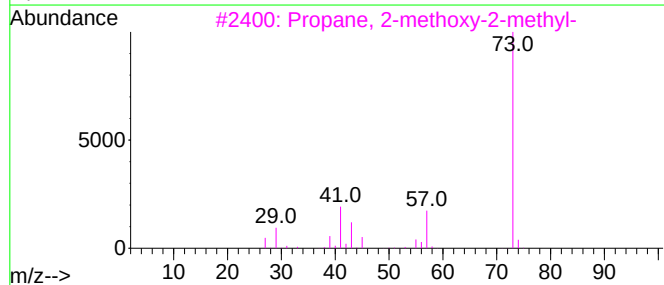
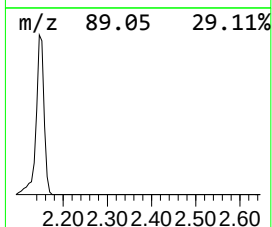
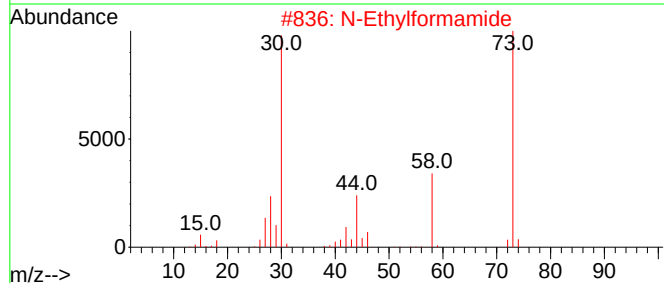
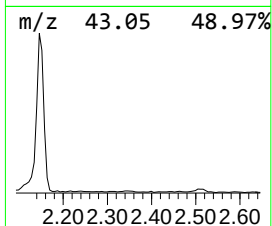
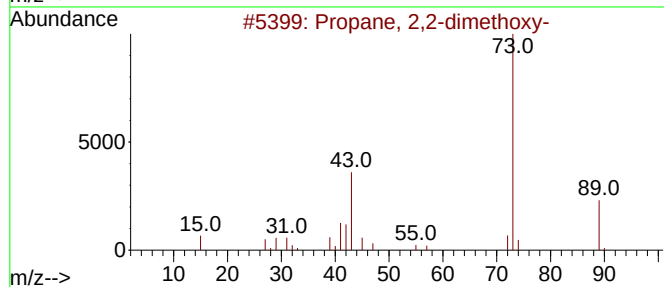
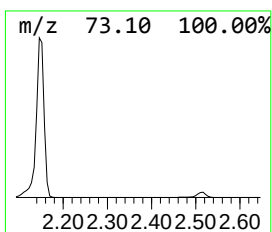
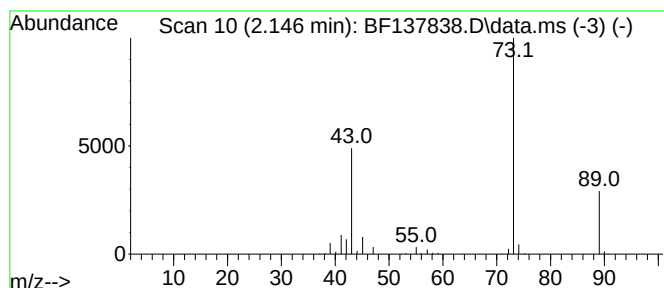
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Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.146	7.27 ng	304092	1,4-Dichlorobenzene-d4	7.046

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	50
2			N-Ethylformamide	73	C3H7NO	000627-45-2	9
3			Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
4			2-Acetamido-N-methylacetamide	130	C5H10N2O2	007606-79-3	9
5			2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9



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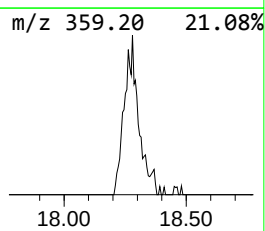
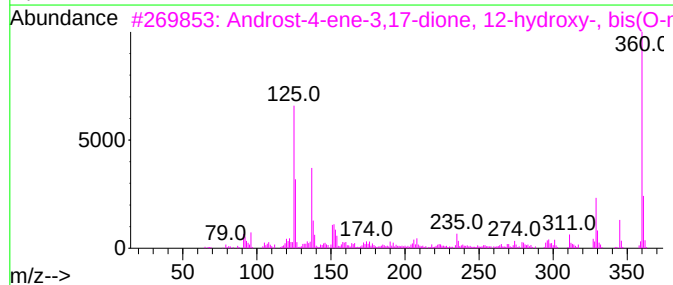
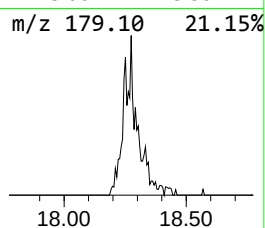
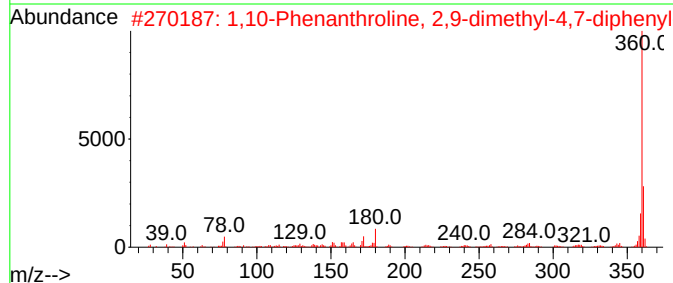
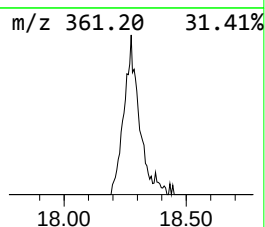
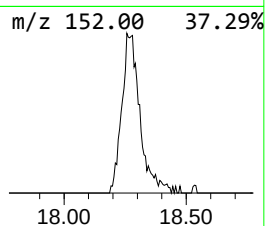
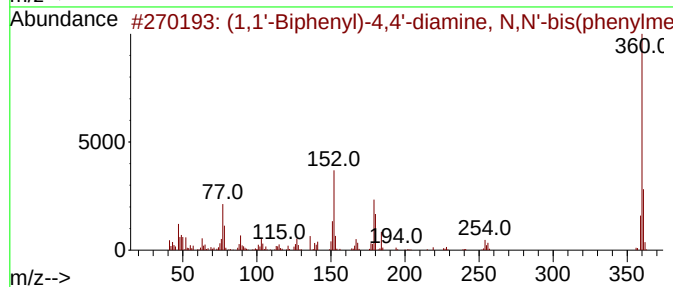
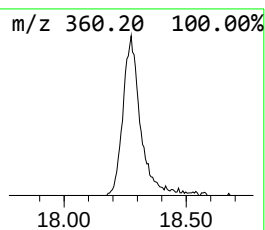
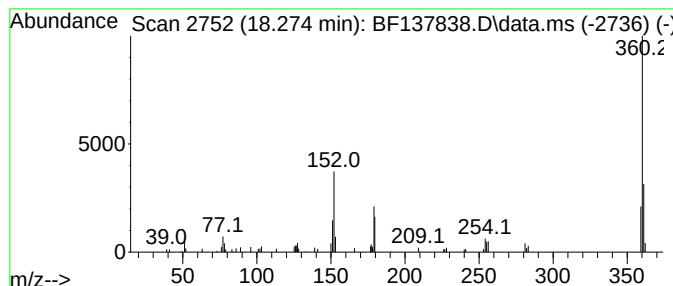
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Peak Number 2 (1,1'-Biphenyl)-4,4'-diamin... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.274	4.04 ng	199803	Perylene-d12	15.804

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(1,1'-Biphenyl)-4,4'-diamine, N,...	360	C26H20N2	006311-48-4	94
2			1,10-Phenanthroline, 2,9-dimethy...	360	C26H20N2	004733-39-5	58
3			Androst-4-ene-3,17-dione, 12-hyd...	360	C21H32N2O3	069688-31-9	50
4			Benzaldehyde, 4-hydroxy-3,5-dime...	360	C18H20N2O6	014414-32-5	50
5			Pentacyclo[18.4.1.1(2,7).1(8,13)...	360	C28H24	101077-60-5	50



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 2,2-di...	2.146	7.3	ng	304092	1	7.046	836023	20.0
(1,1'-Biphenyl)...	18.274	4.0	ng	199803	6	15.804	987908	20.0