

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF073121\
 Data File : BF124893.D
 Acq On : 31 Jul 2021 15:24
 Operator : JU/CG
 Sample : M3235-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-462-COMP

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: BF124893.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.075	505	508	512	rBB	2744	2641	1.71%	0.287%
2	5.475	571	576	579	rBB	132903	131571	85.15%	14.288%
3	6.487	743	748	750	rBV	133493	136648	88.44%	14.839%
4	6.857	808	811	813	rBB	32857	25998	16.83%	2.823%
5	7.416	902	906	909	rBB	94900	87939	56.91%	9.550%
6	8.034	1009	1011	1016	rBB	15905	12191	7.89%	1.324%
7	8.139	1025	1029	1032	rBB	44785	35968	23.28%	3.906%
8	9.210	1207	1211	1214	rBV	190210	154516	100.00%	16.780%
9	9.892	1324	1327	1329	rBB	50958	39420	25.51%	4.281%
10	10.680	1457	1461	1464	rBB	110804	88454	57.25%	9.606%
11	11.375	1576	1579	1582	rBV	49104	37752	24.43%	4.100%
12	12.963	1845	1849	1852	rBB	144373	124721	80.72%	13.544%
13	14.010	2024	2027	2031	rBB	22930	20831	13.48%	2.262%
14	15.480	2272	2277	2281	rBB	19024	22202	14.37%	2.411%

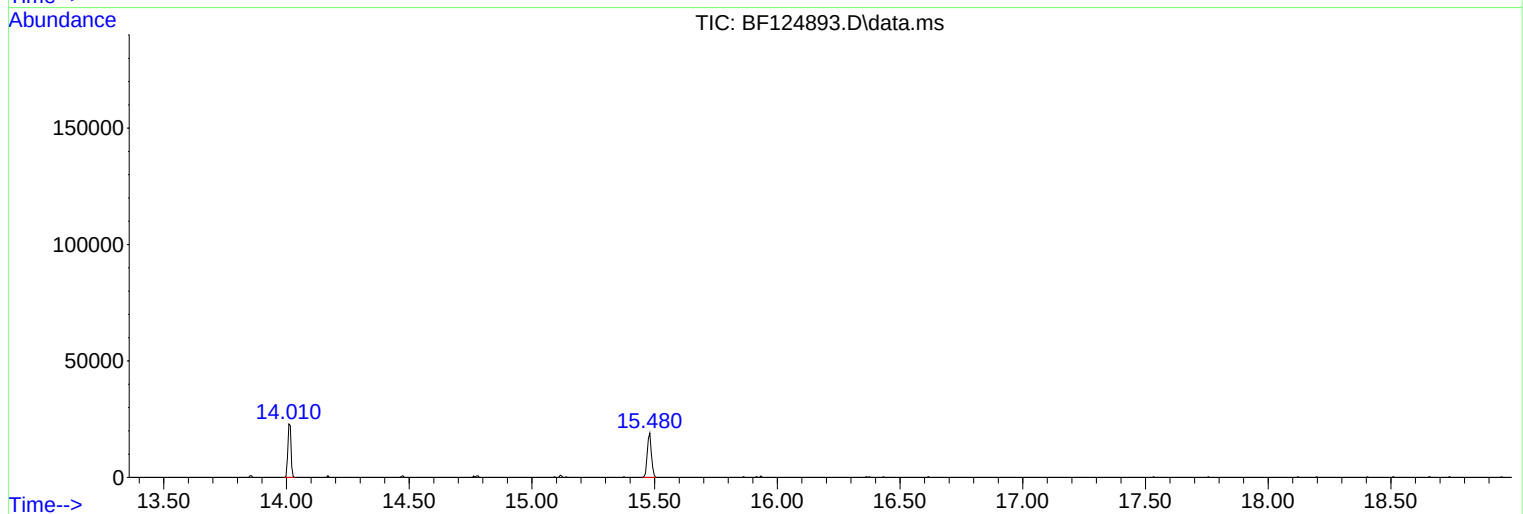
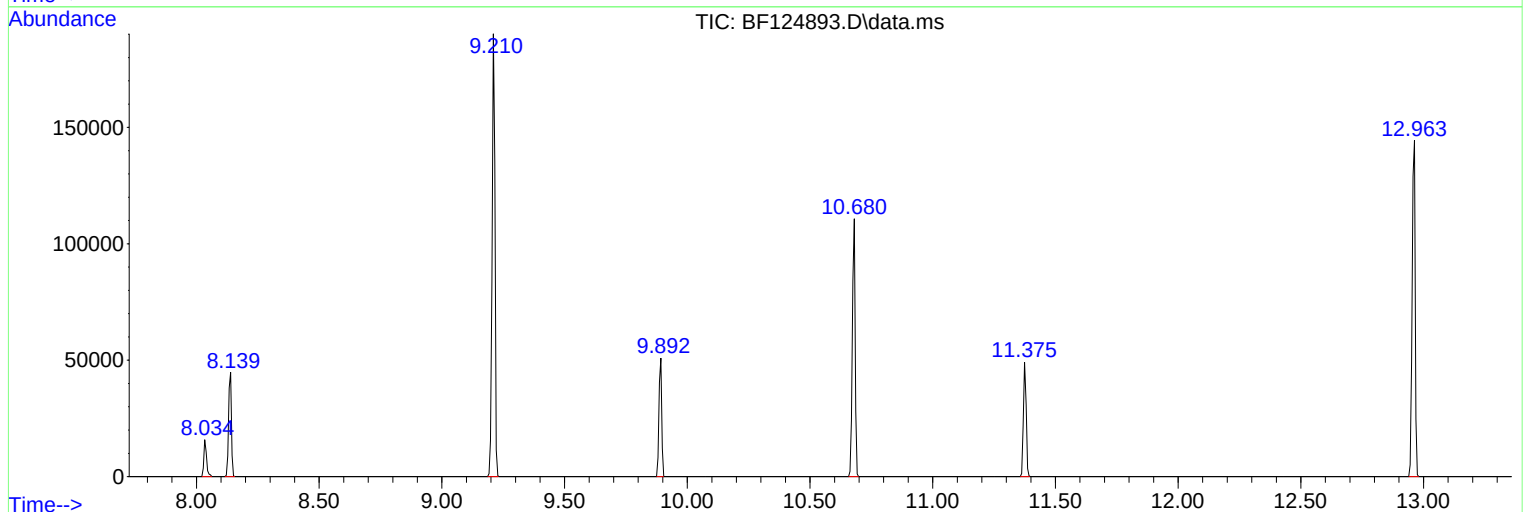
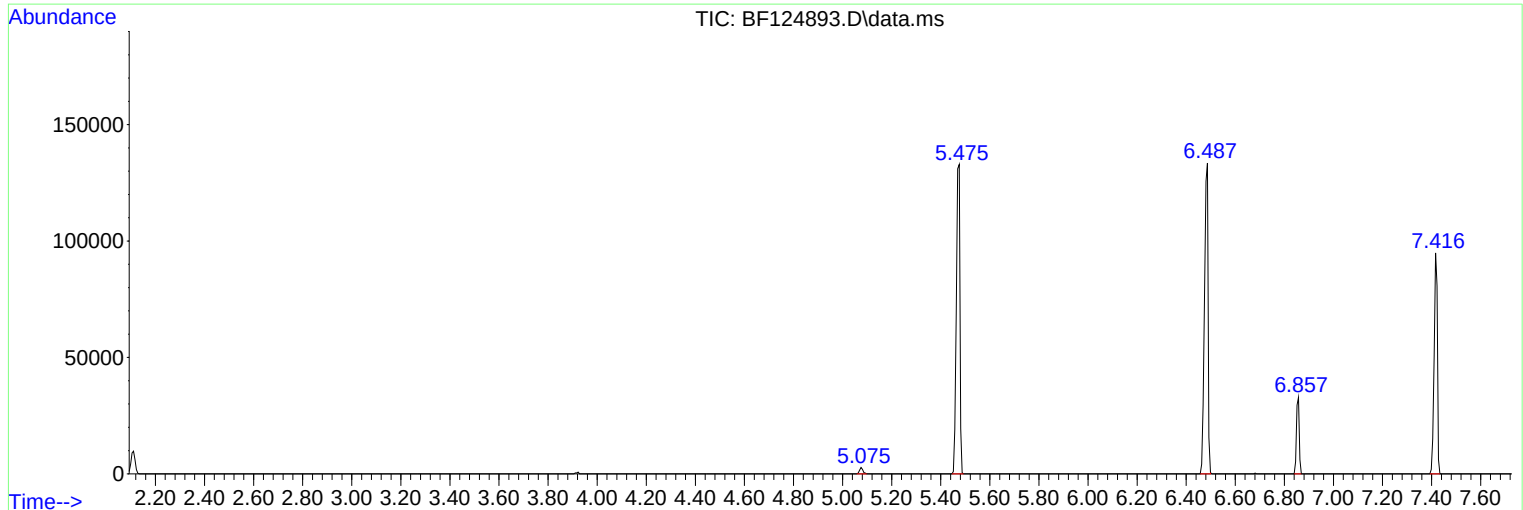
Sum of corrected areas: 920852

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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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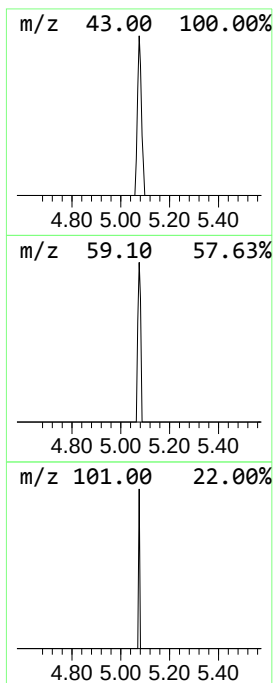
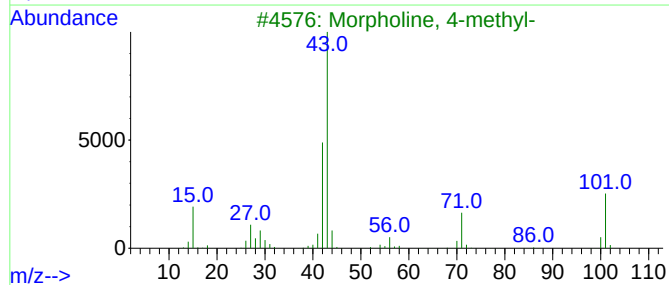
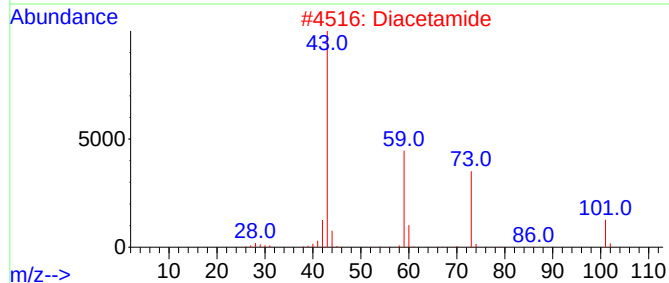
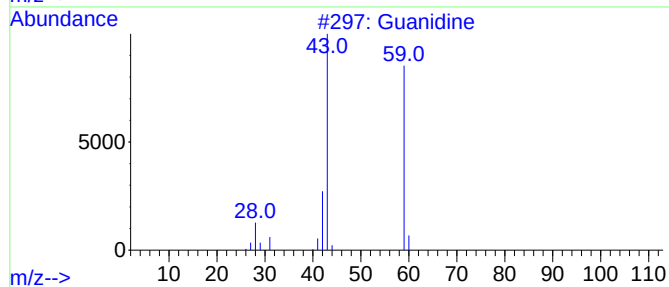
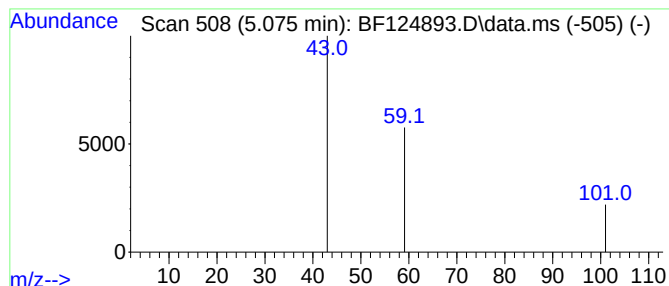
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Peak Number 1 unknown5.075 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.075	2.03 ng	2641	1,4-Dichlorobenzene-d4	6.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Guanidine	59	CH5N3	000113-00-8	4
2			Diacetamide	101	C4H7NO2	000625-77-4	4
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	4
4			Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	4
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	4



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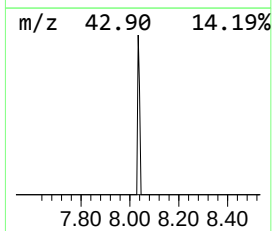
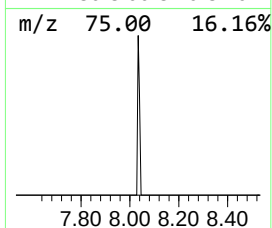
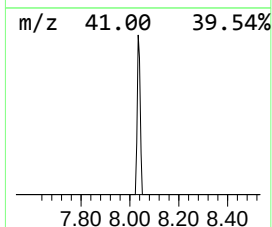
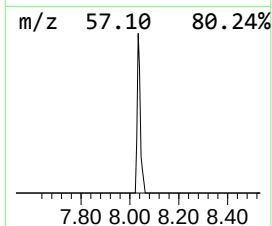
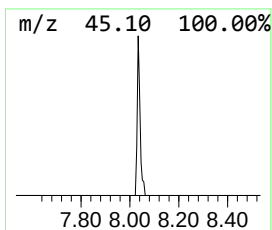
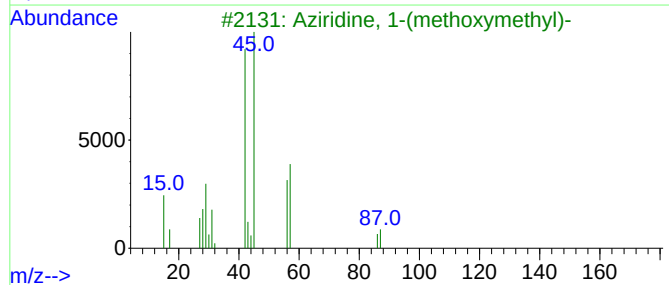
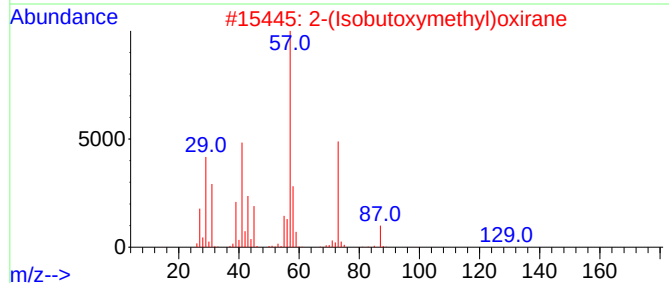
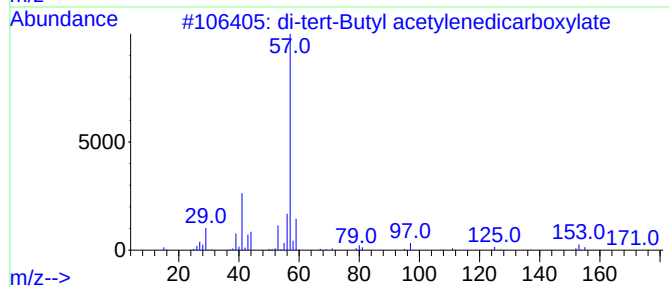
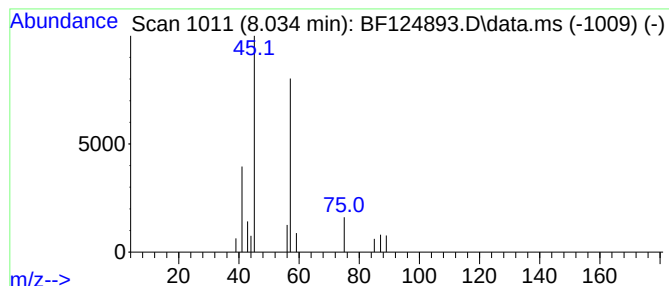
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Peak Number 2 unknown8.034 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.034	6.78 ng	12191	Naphthalene-d8	8.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			di-tert-Butyl acetylenedicarboxy...	226	C12H18O4	066086-33-7	9
2			2-(Isobutoxymethyl)oxirane	130	C7H14O2	003814-55-9	9
3			Aziridine, 1-(methoxymethyl)-	87	C4H9NO	001497-83-2	7
4			4-Amino-1-butanol	89	C4H11NO	013325-10-5	5
5			o-tert-Butylhydroxylamine	89	C4H11NO	1000322-43-1	5



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown5.075	5.075	2.0	ng	2641	1	6.857	25998	20.0
unknown8.034	8.034	6.8	ng	12191	2	8.139	35968	20.0