

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072321\
 Data File : BF124717.D
 Acq On : 23 Jul 2021 17:19
 Operator : JU/CG
 Sample : PB137842BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB137842BL

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: BF124717.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.134	515	518	523	rBB	3332	4723	1.34%	0.248%
2	5.498	575	580	583	rBB	214535	250943	71.38%	13.158%
3	6.498	743	750	752	rBB	215356	249973	71.10%	13.107%
4	6.869	810	813	816	rBB	56881	43065	12.25%	2.258%
5	7.433	903	909	911	rBB	151015	163376	46.47%	8.567%
6	8.151	1026	1031	1034	rBB	64875	60042	17.08%	3.148%
7	9.227	1208	1214	1217	rBB	313274	332913	94.70%	17.457%
8	9.904	1324	1329	1332	rBB	85245	71946	20.46%	3.773%
9	10.698	1458	1464	1467	rBB	218329	204540	58.18%	10.725%
10	11.392	1577	1582	1585	rBB	89409	71980	20.47%	3.774%
11	12.986	1848	1853	1856	rVB	363385	351557	100.00%	18.434%
12	14.033	2027	2031	2034	rBB	62274	55394	15.76%	2.905%
13	15.504	2276	2281	2285	rBB	38839	46648	13.27%	2.446%

Sum of corrected areas: 1907100

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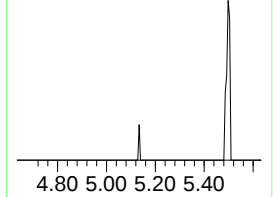
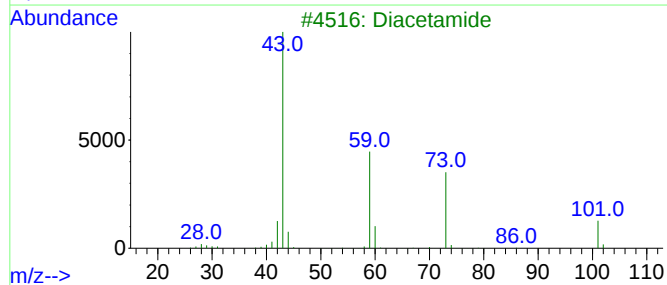
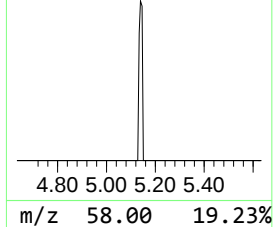
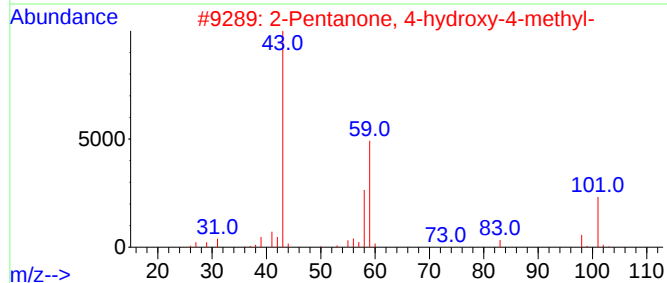
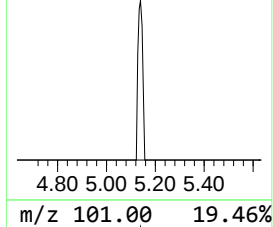
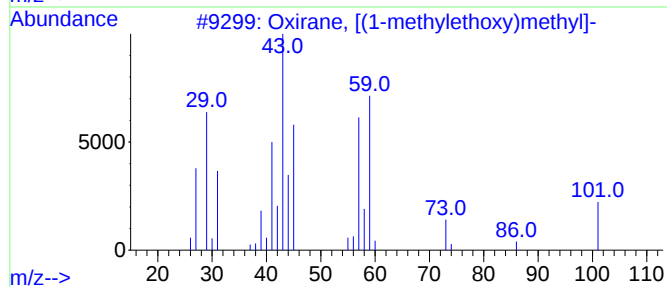
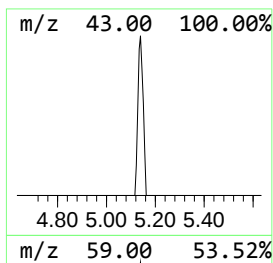
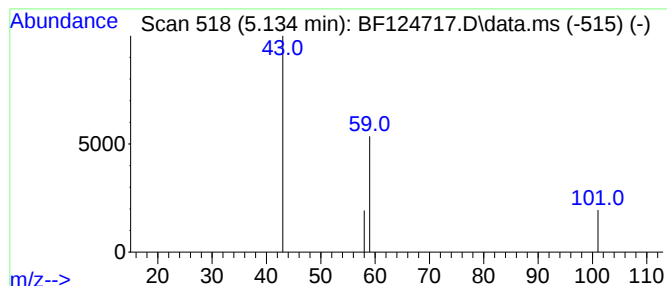
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown5.134 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.134	2.19 ng	4723	1,4-Dichlorobenzene-d4	6.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	9		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	9		
3	Diacetamide	101	C4H7NO2	000625-77-4	4		
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	4		
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	3		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown5.134	5.134	2.2	ng	4723	1	6.869	43065	20.0