

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080321\
 Data File : BF124922.D
 Acq On : 3 Aug 2021 18:50
 Operator : JU/CG
 Sample : PB138173BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB138173BL

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: BF124922.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	515	rBB	12916	17082	6.24%	1.163%
2	5.469	571	575	578	rBV	185613	191152	69.77%	13.020%
3	6.475	741	746	749	rBV	174401	191316	69.83%	13.031%
4	6.845	806	809	812	rBB	47617	34847	12.72%	2.373%
5	7.410	900	905	908	rBB	129837	125018	45.63%	8.515%
6	8.128	1023	1027	1030	rBB	59444	47096	17.19%	3.208%
7	9.204	1205	1210	1213	rBB	282544	245198	89.50%	16.701%
8	9.880	1322	1325	1328	rBB	68226	54135	19.76%	3.687%
9	10.674	1455	1460	1463	rBV	170786	147382	53.80%	10.038%
10	11.368	1574	1578	1581	rBB	72930	57286	20.91%	3.902%
11	12.957	1843	1848	1851	rBV	314168	273969	100.00%	18.661%
12	13.868	2000	2003	2006	rBB	8313	7620	2.78%	0.519%
13	14.004	2023	2026	2029	rBB	53093	43100	15.73%	2.936%
14	15.468	2271	2275	2279	rBB	31322	32973	12.04%	2.246%

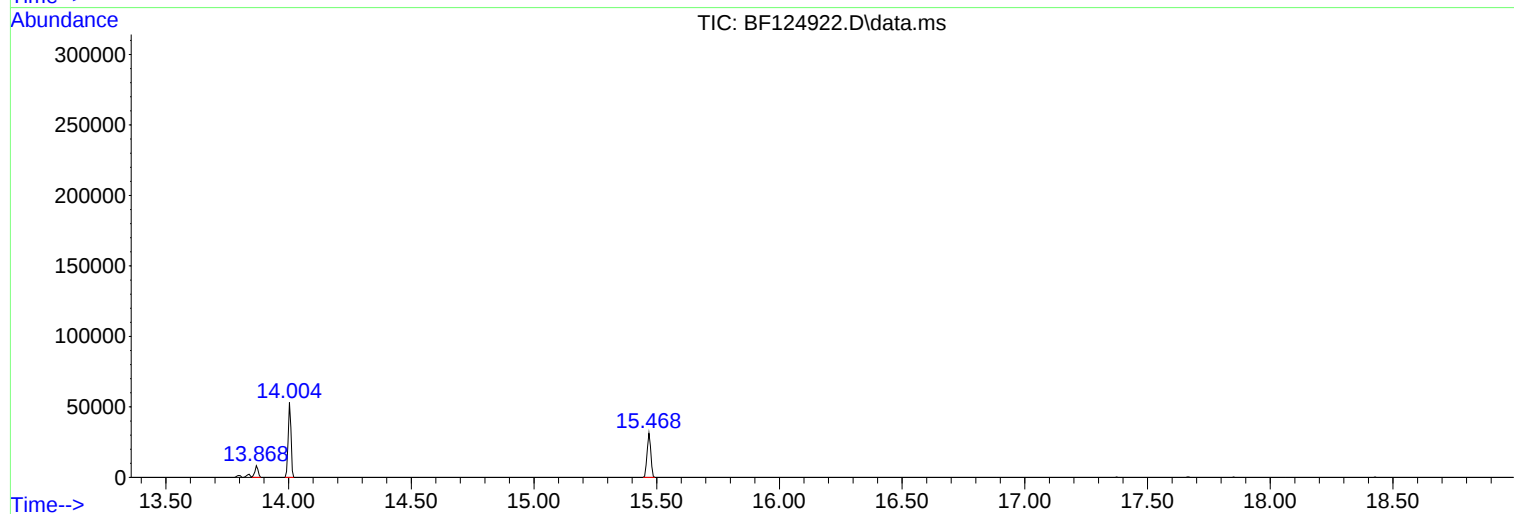
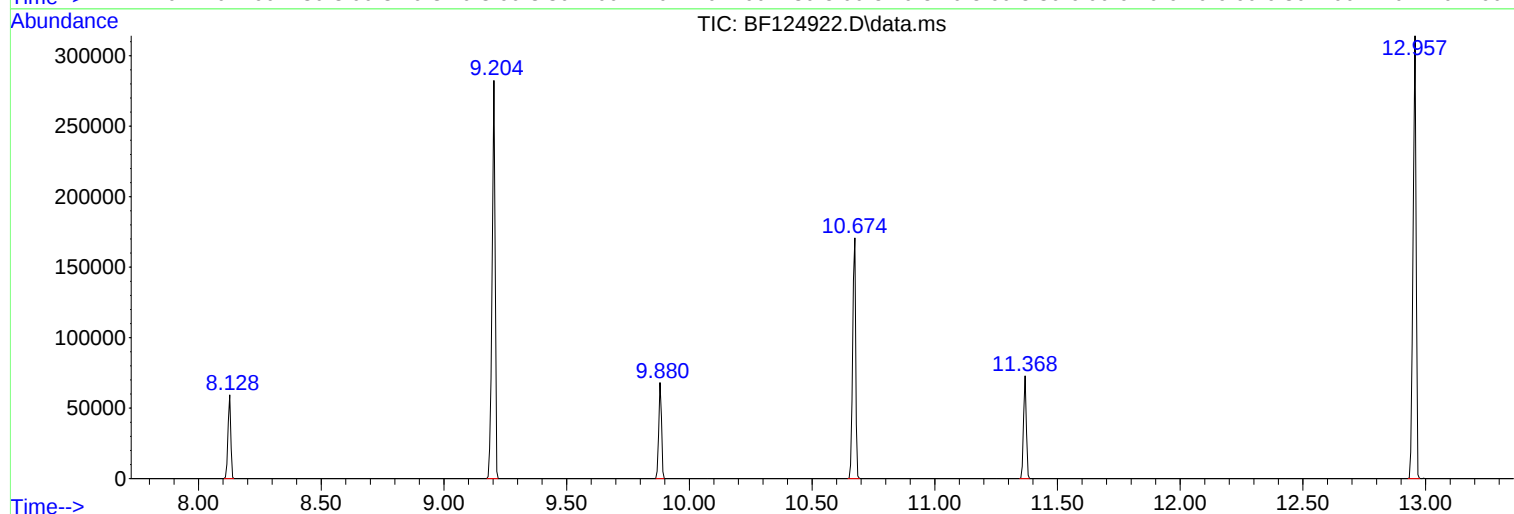
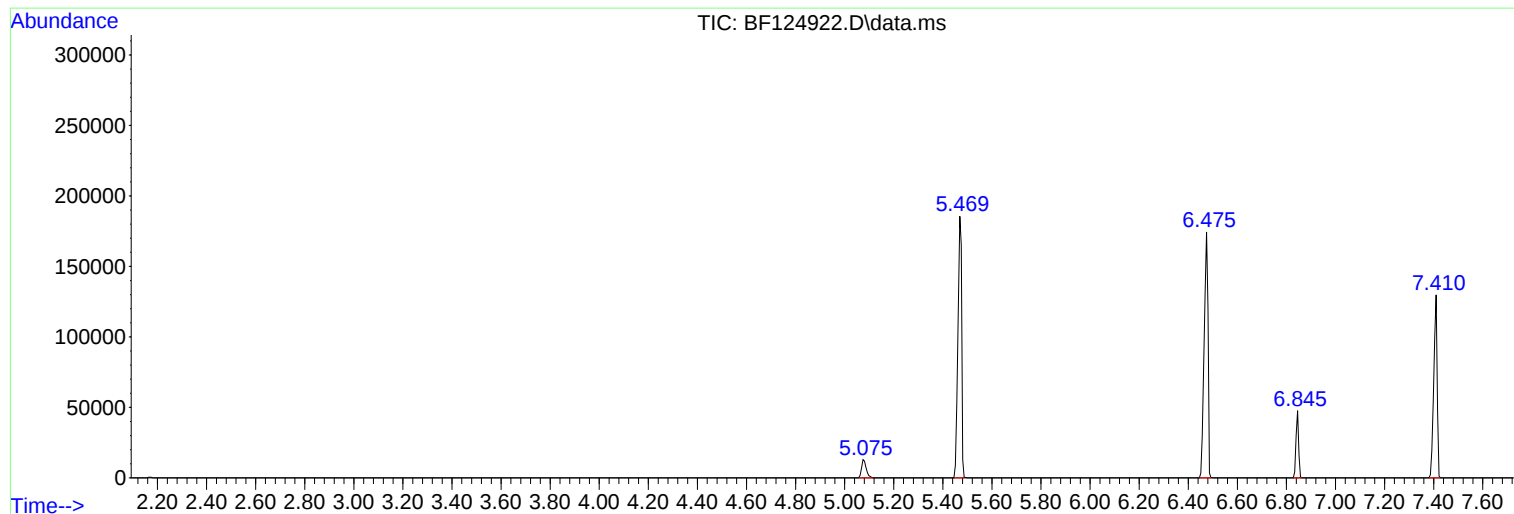
Sum of corrected areas: 1468174

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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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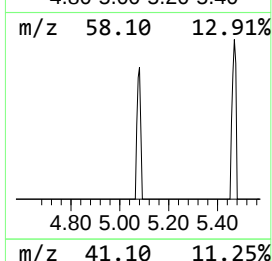
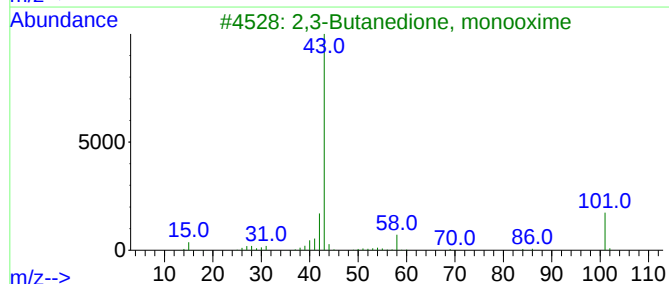
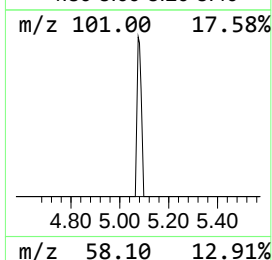
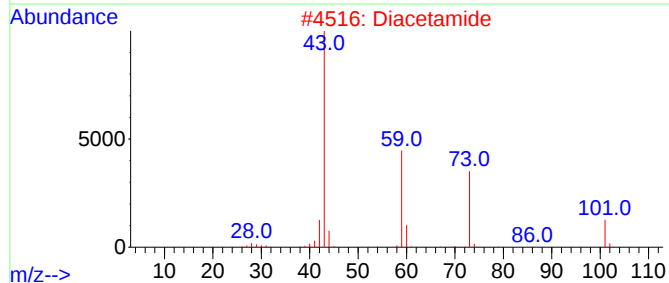
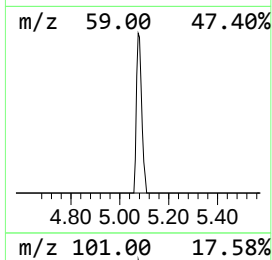
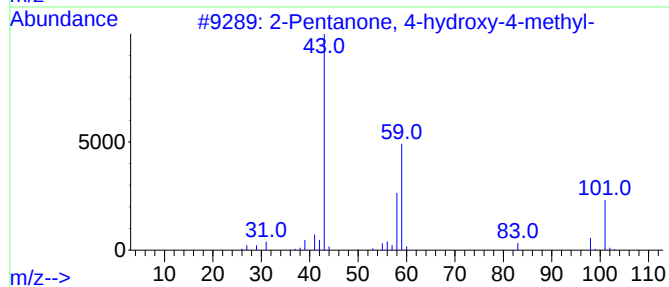
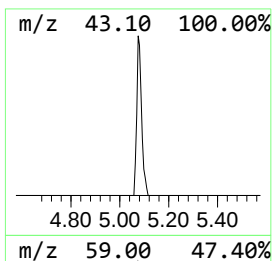
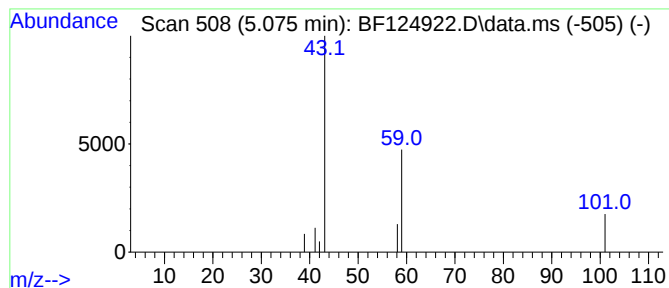
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Peak Number 1 unknown5.075 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.075	9.80 ng	17082	1,4-Dichlorobenzene-d4	6.845

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	9		
2	Diacetamide	101	C4H7NO2	000625-77-4	7		
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	7		
4	N-(n-Propyl)acetamide	101	C5H11NO	005331-48-6	4		
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	4		



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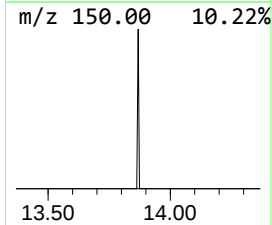
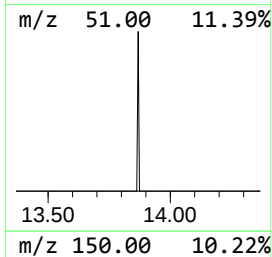
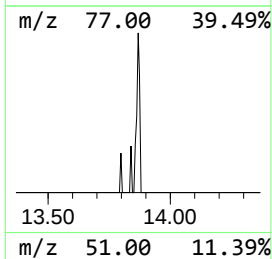
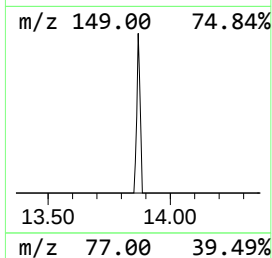
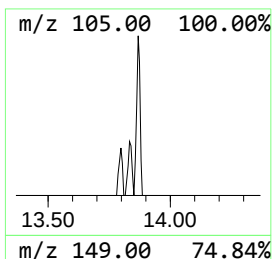
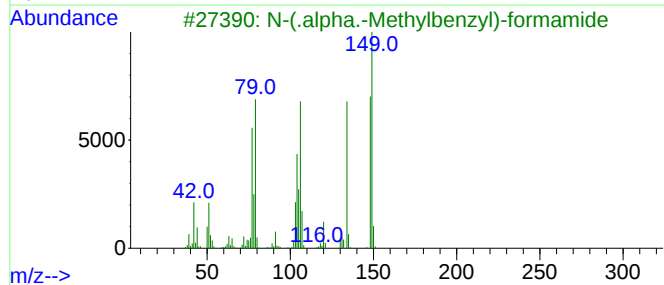
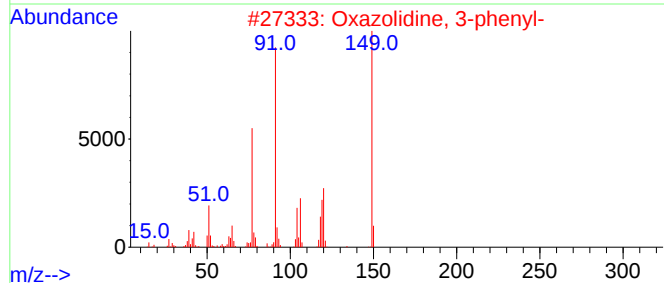
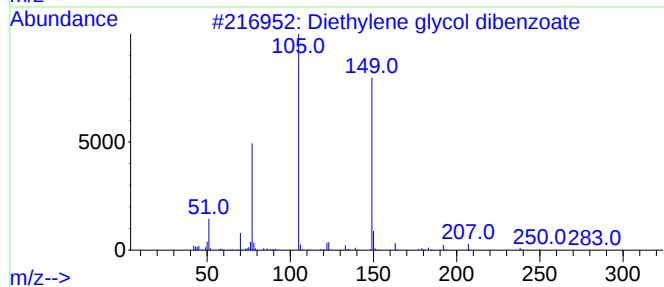
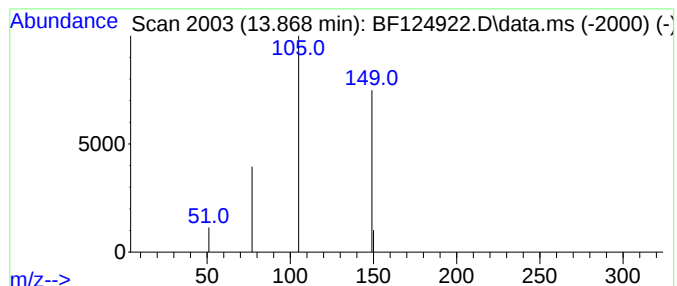
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Peak Number 2 Diethylene glycol dibenzoate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.868	3.54 ng	7620	Chrysene-d12	14.004	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	74
2	Oxazolidine, 3-phenyl-	149	C9H11NO	020503-92-8	74
3	N-(.alpha.-Methylbenzyl)-formamide	149	C9H11NO	006948-01-2	74
4	2-(2-(2-Methoxyethoxy)ethoxy)eth...	268	C14H20O5	1000366-99-1	64
5	2,2'-(Ethane-1,2-diylbis(oxy))bi...	358	C20H22O6	1000366-99-4	64



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown5.075	5.075	9.8	ng	17082	1	6.845	34847	20.0
Diethylene glyc...	13.868	3.5	ng	7620	5	14.004	43100	20.0