

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062918\
 Data File : BG035528.D
 Acq On : 30 Jun 2018 5:47
 Operator : JU/SJ
 Sample : J3774-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-1(4-6)

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_G\METHODS\8270-BG060718.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.171	315	320	327	rBV	39478	63630	3.14%	0.631%
2	5.641	392	400	410	rBV	403474	651625	32.14%	6.463%
3	7.221	660	669	685	rBV	435419	766033	37.78%	7.598%
4	7.597	726	733	744	rBV	405661	692672	34.16%	6.870%
5	8.061	806	812	820	rBV	85779	147390	7.27%	1.462%
6	8.384	861	867	875	rBV	278368	482714	23.81%	4.788%
7	9.224	1000	1010	1019	rBV	261542	455872	22.48%	4.522%
8	10.863	1282	1289	1296	rBV	115727	212672	10.49%	2.109%
9	13.307	1698	1705	1714	rBV	726377	1106216	54.56%	10.972%
10	14.129	1840	1845	1851	rBV2	60911	90332	4.46%	0.896%
11	14.682	1933	1939	1947	rBV2	229195	369514	18.22%	3.665%
12	16.168	2185	2192	2202	rBV	725583	1106766	54.58%	10.978%
13	17.425	2396	2406	2414	rBV	365565	550098	27.13%	5.456%
14	18.306	2552	2556	2561	rBV3	35737	48547	2.39%	0.482%
15	20.027	2843	2849	2855	rBV	1573969	2027651	100.00%	20.112%
16	21.326	3066	3070	3075	rBV3	24999	41214	2.03%	0.409%
17	21.690	3125	3132	3140	rVB	395623	627150	30.93%	6.221%
18	23.370	3412	3418	3425	rBV3	27476	53382	2.63%	0.529%
19	24.915	3670	3681	3694	rVB2	202545	588412	29.02%	5.836%

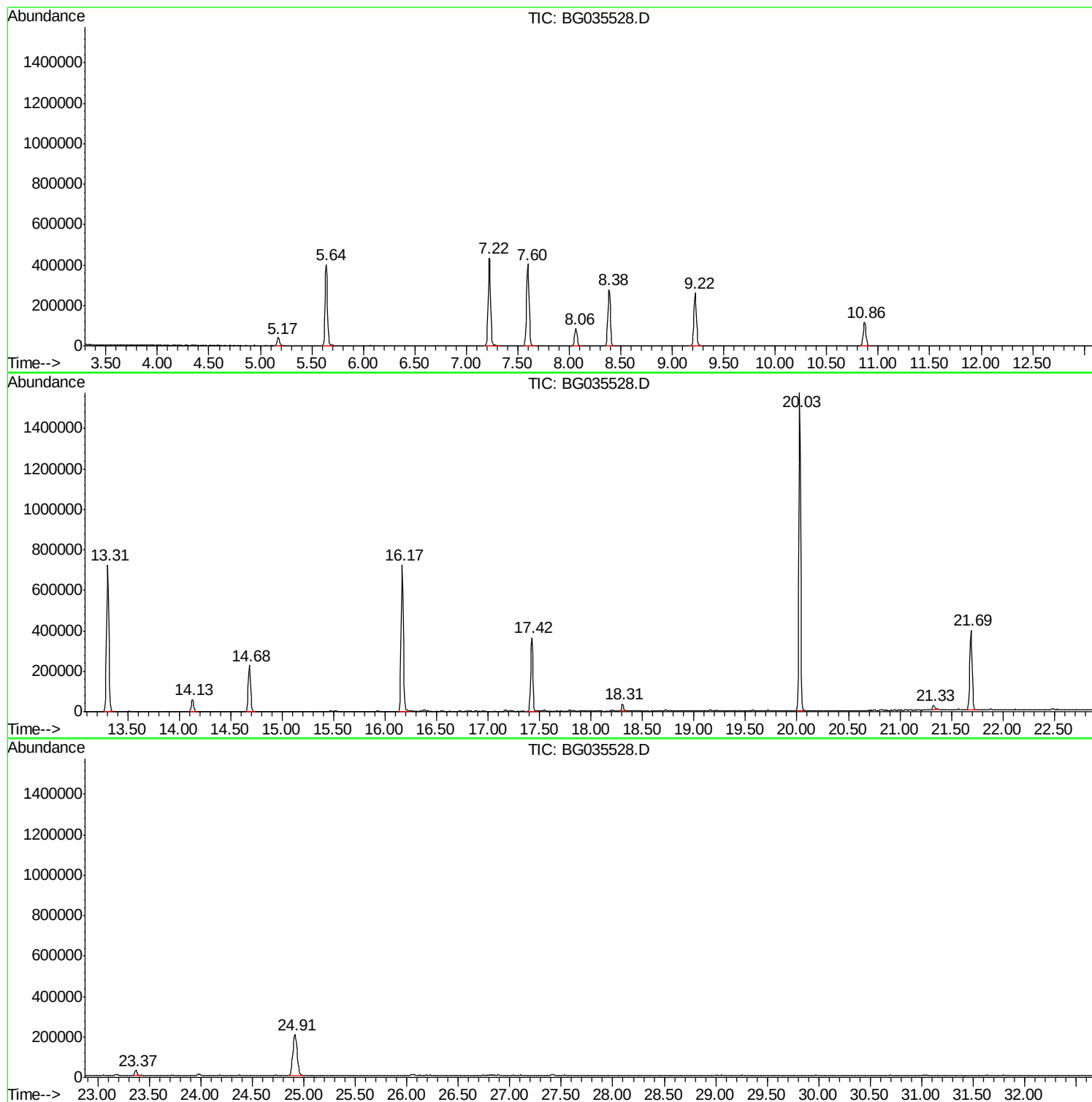
Sum of corrected areas: 10081890

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG060718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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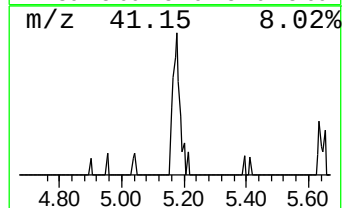
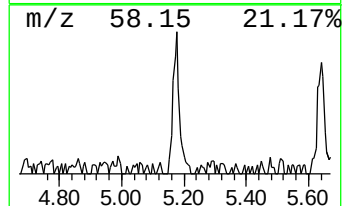
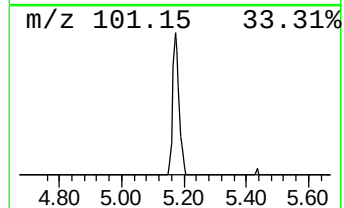
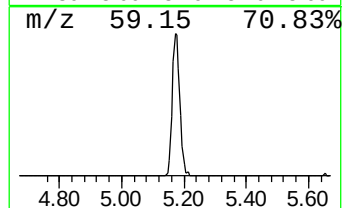
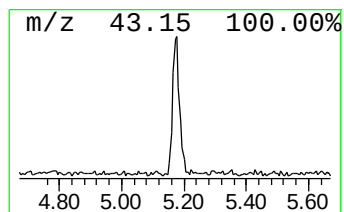
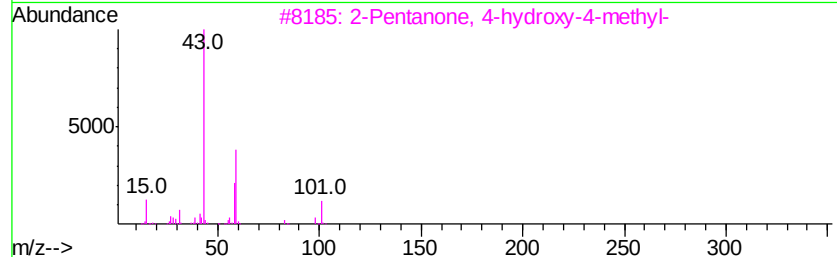
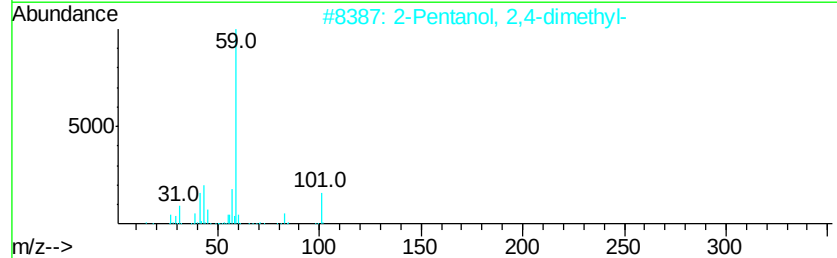
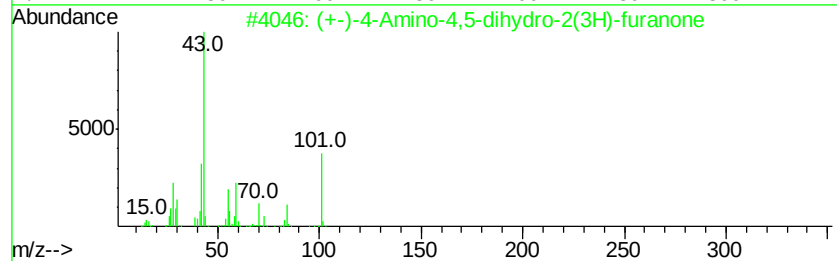
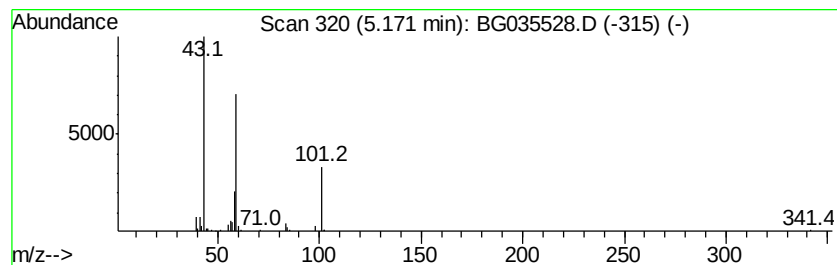
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	8.63 ng	63630	1,4-Dichlorobenzene-d4	8.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-	4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	47	
2	2-Pentanol, 2,4-dimethyl-		116	C7H16O	000625-06-9	47	
3	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	36	
4	4-Ethyl-3-octanol		158	C10H22O	019781-26-1	25	
5	2-Hexanol, 2-methyl-		116	C7H16O	000625-23-0	25	



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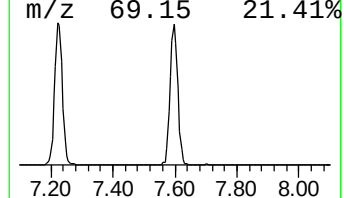
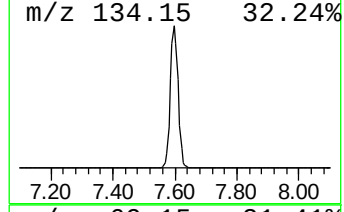
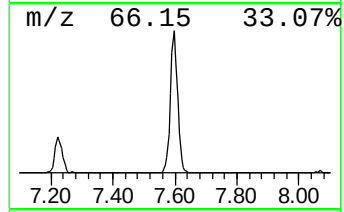
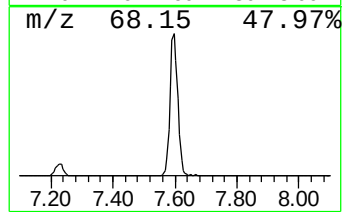
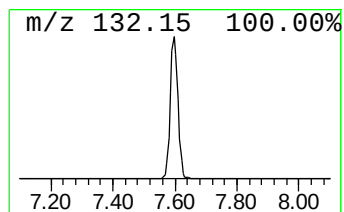
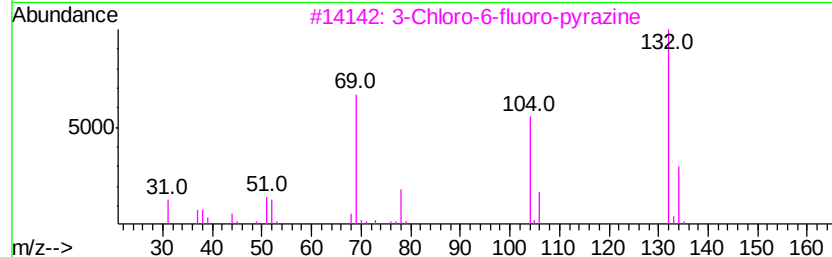
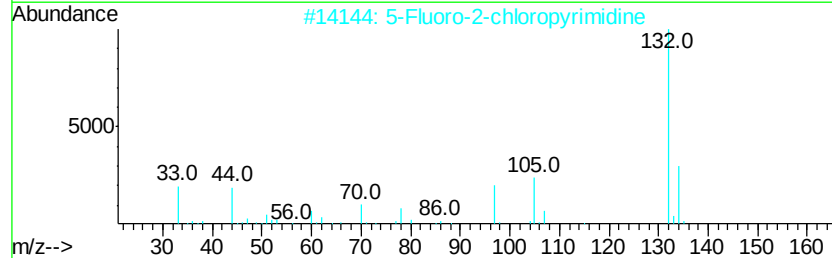
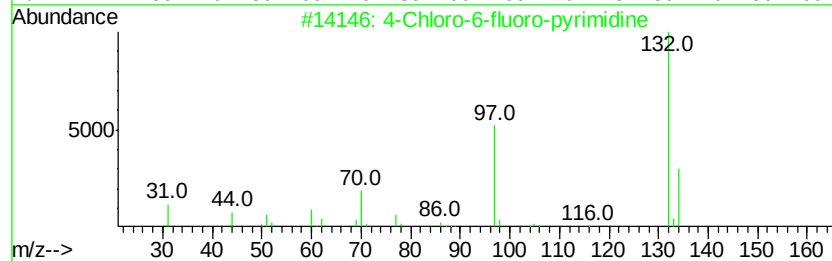
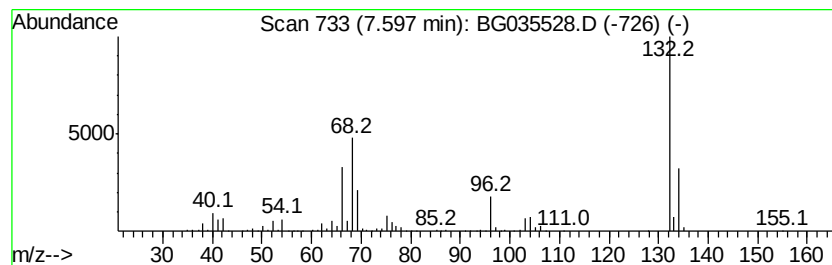
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Peak Number 2 unknown7.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	93.99 ng	692672	1,4-Dichlorobenzene-d4	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
4		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		1H-Indazole, 3-methyl-	132	C8H8N2	1000316-00-2	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	8.6	ng	63630	1	8.07	147390	20.0
unknown7.60	7.60	94.0	ng	692672	1	8.07	147390	20.0