

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102918\
 Data File : BG037617.D
 Acq On : 29 Oct 2018 11:18
 Operator : JU/SJ
 Sample : PB114233BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114233BL

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-BG101818.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	5.180	315	322	333	rVB	33014	60050	1.80%	0.366%
2	5.644	394	401	413	rBV	750502	1237594	37.07%	7.537%
3	7.242	666	673	687	rBV	777048	1401845	41.99%	8.537%
4	7.630	731	739	748	rBV	716282	1293713	38.75%	7.879%
5	8.100	813	819	827	rBV	146590	261568	7.83%	1.593%
6	8.423	866	874	883	rBV	528978	965756	28.93%	5.882%
7	9.275	1011	1019	1031	rBV	443639	831764	24.91%	5.066%
8	10.920	1292	1299	1309	rVB	210262	398669	11.94%	2.428%
9	13.359	1706	1714	1724	rBV	1346600	2144394	64.23%	13.060%
10	14.734	1939	1948	1956	rBV2	366193	614647	18.41%	3.743%
11	16.220	2194	2201	2213	rBV2	935824	1513059	45.32%	9.215%
12	17.483	2409	2416	2426	rBV2	507767	819398	24.54%	4.990%
13	20.080	2852	2858	2872	rBV	2503317	3338580	100.00%	20.333%
14	21.767	3138	3145	3157	rBV	475076	856906	25.67%	5.219%
15	25.104	3701	3713	3726	rBV3	186923	681973	20.43%	4.153%

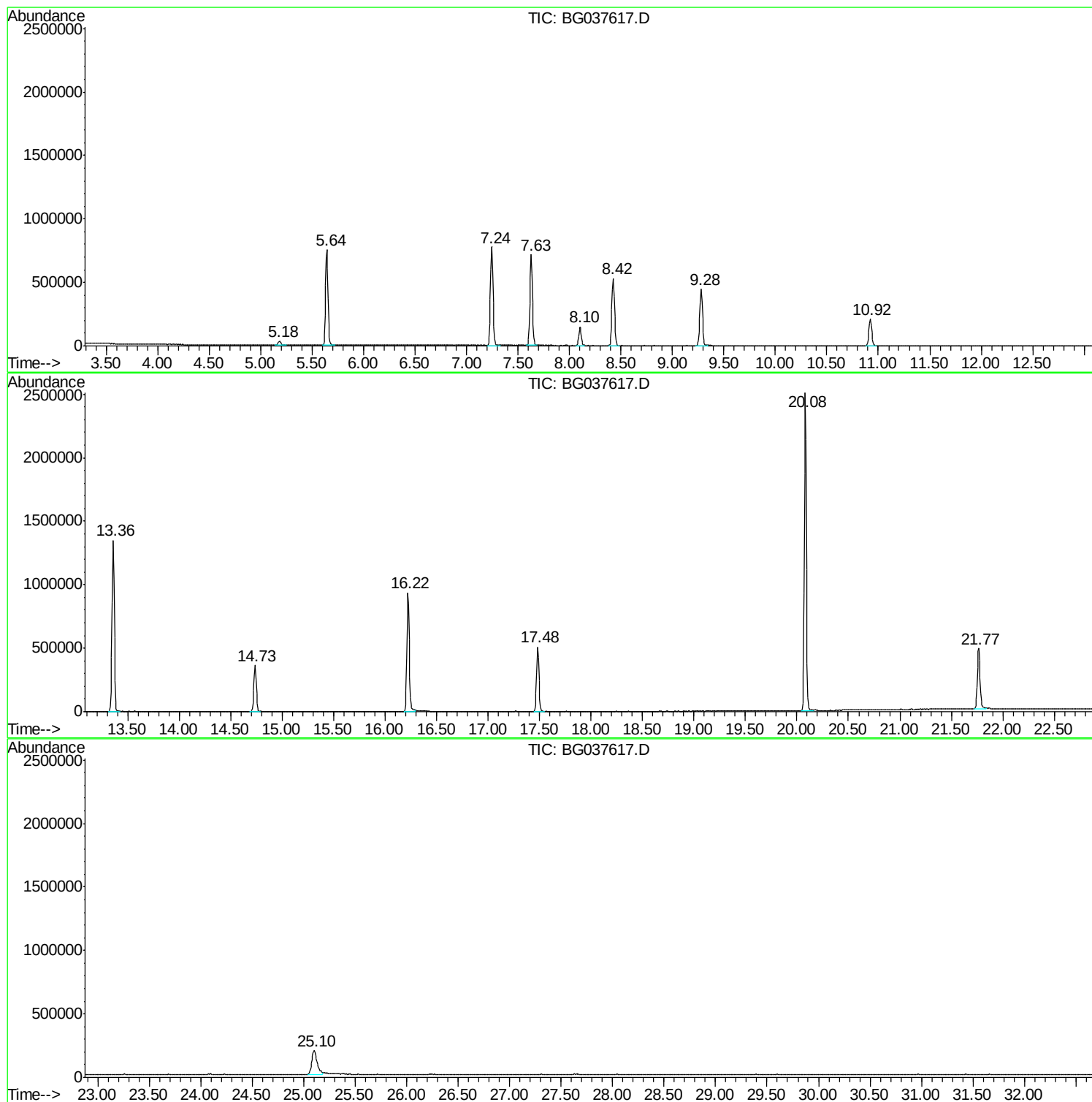
Sum of corrected areas: 16419916

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.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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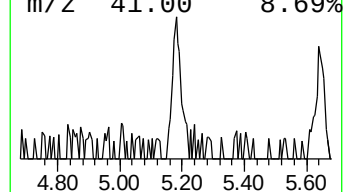
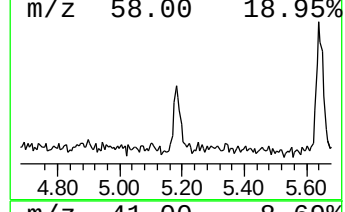
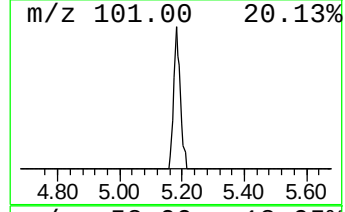
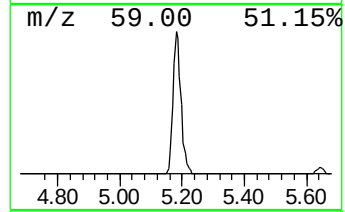
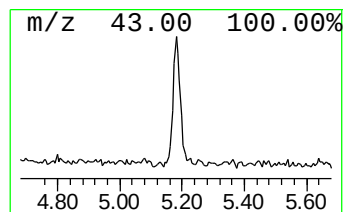
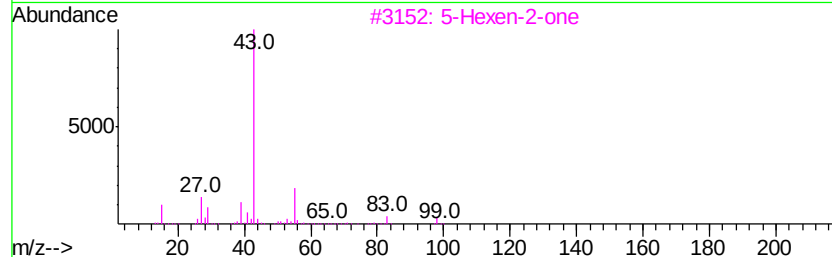
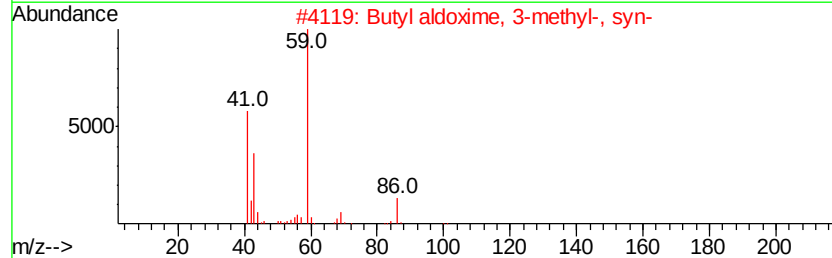
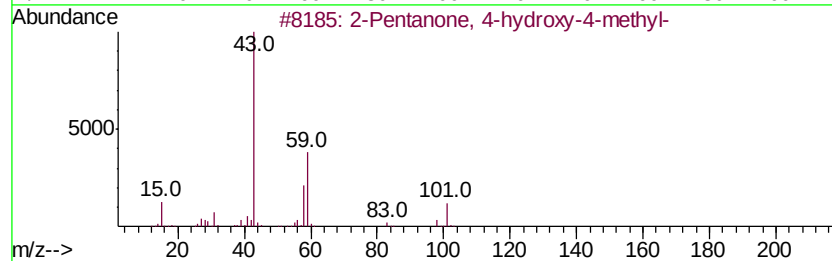
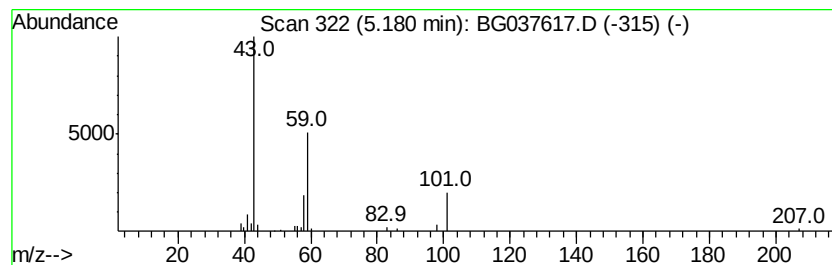
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	4.59 ng	60050	1,4-Dichlorobenzene-d4	8.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	12
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			3-Octanol	130	C8H18O	000589-98-0	9



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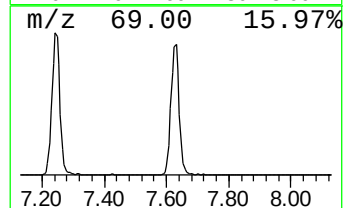
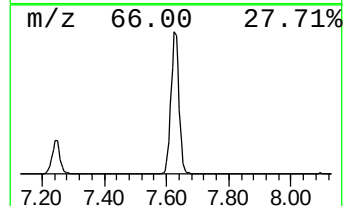
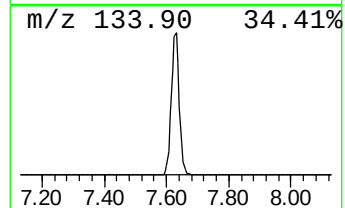
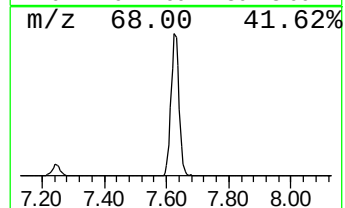
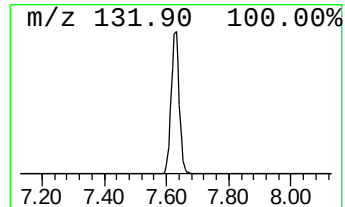
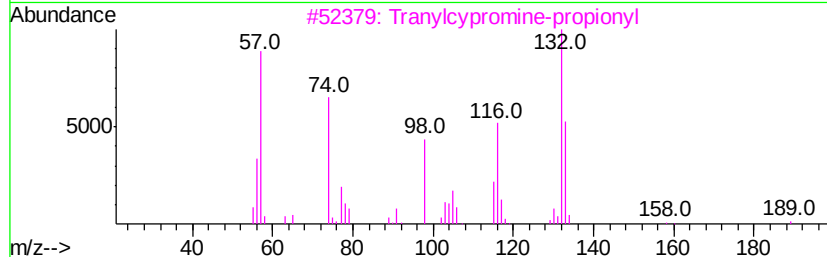
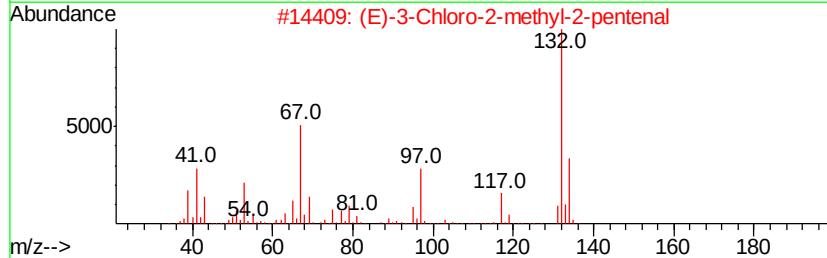
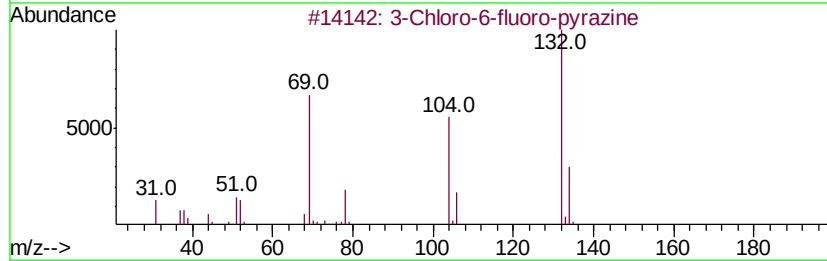
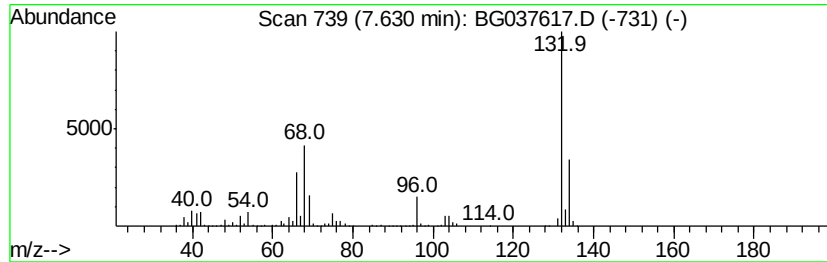
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Peak Number 2 unknown7.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	98.92 ng	1293710	1,4-Dichlorobenzene-d4	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		Amphetaminil	250	C17H18N2	017590-01-1	10
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	4.6	ng	60050	1	8.10	261568	20.0
unknown7.63	7.63	98.9	ng	1293710	1	8.10	261568	20.0