

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110218\
 Data File : BG037793.D
 Acq On : 3 Nov 2018 16:15
 Operator : JU/SJ
 Sample : PB114415BL
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114415BL

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	5.179	316	322	328	rBV2	39849	67106	1.77%	0.337%
2	5.637	393	400	413	rBV	959083	1577976	41.62%	7.920%
3	7.241	664	673	685	rBV	1001041	1795327	47.36%	9.010%
4	7.623	730	738	747	rBV	951291	1693106	44.66%	8.497%
5	8.093	811	818	827	rVB	194601	349759	9.23%	1.755%
6	8.422	865	874	886	rBV	690991	1284745	33.89%	6.448%
7	9.274	1011	1019	1034	rBV	594634	1121990	29.60%	5.631%
8	10.919	1289	1299	1306	rBV	289840	549773	14.50%	2.759%
9	13.352	1705	1713	1725	rBV	1769039	2856788	75.36%	14.338%
10	14.732	1940	1948	1955	rBV2	481580	818945	21.60%	4.110%
11	16.219	2193	2201	2215	rBV	892584	1517378	40.03%	7.615%
12	17.476	2407	2415	2426	rBV	613876	948684	25.02%	4.761%
13	20.079	2851	2858	2875	rBV	2669480	3791045	100.00%	19.026%
14	21.759	3137	3144	3162	rBV	459753	953464	25.15%	4.785%
15	25.097	3700	3712	3722	rBV2	163343	599040	15.80%	3.006%

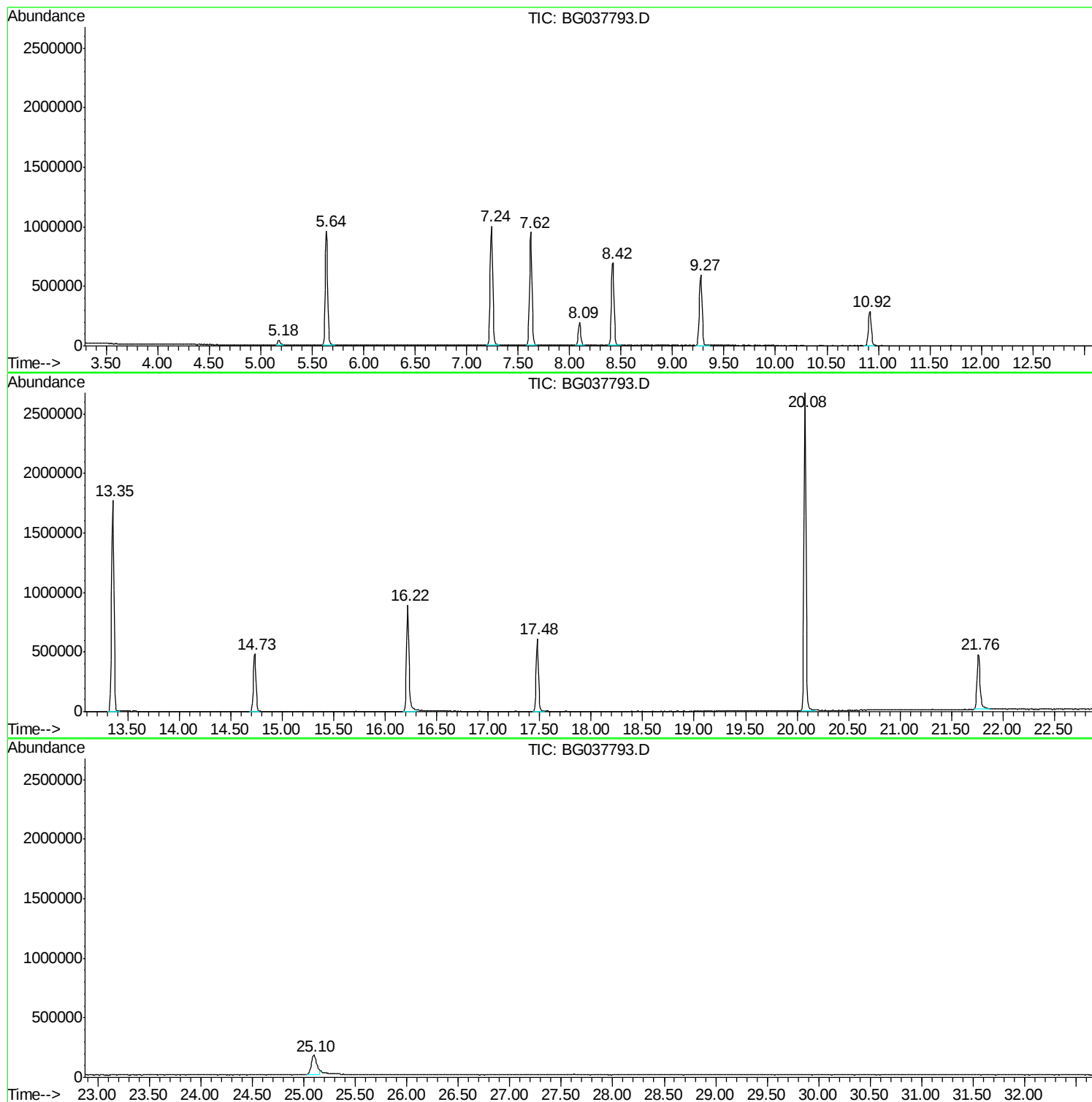
Sum of corrected areas: 19925126

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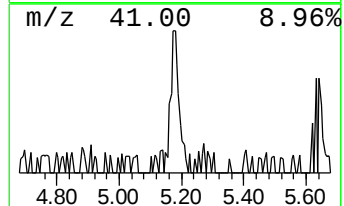
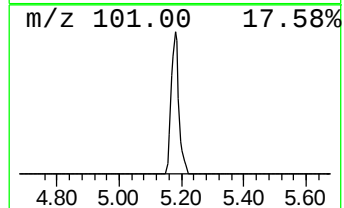
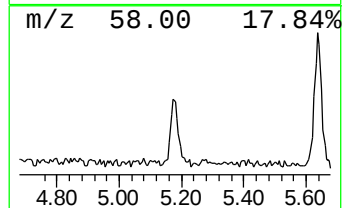
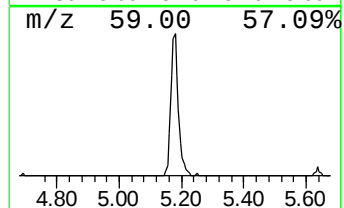
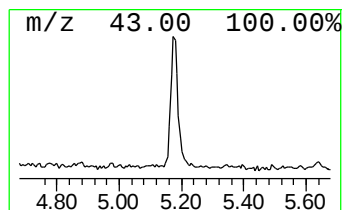
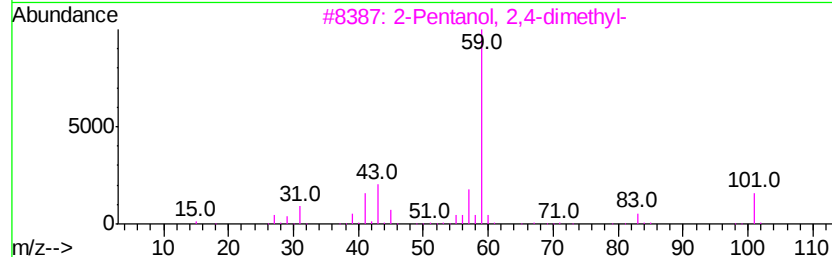
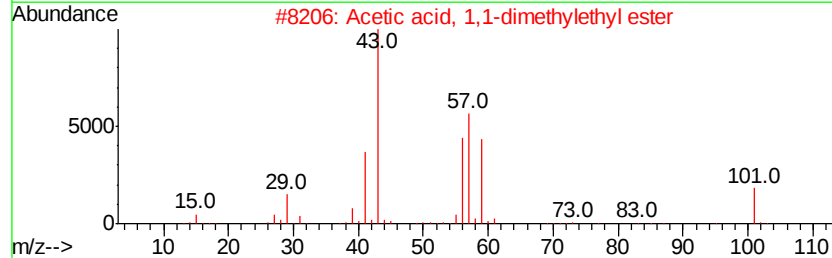
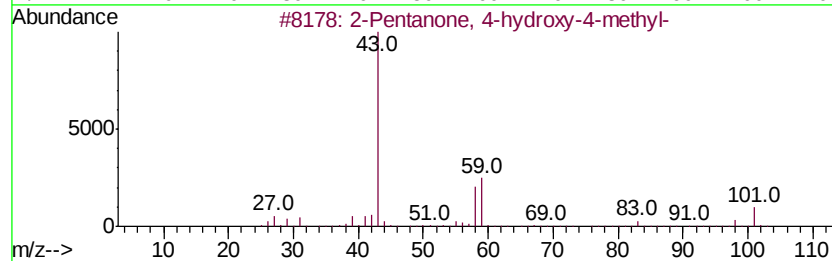
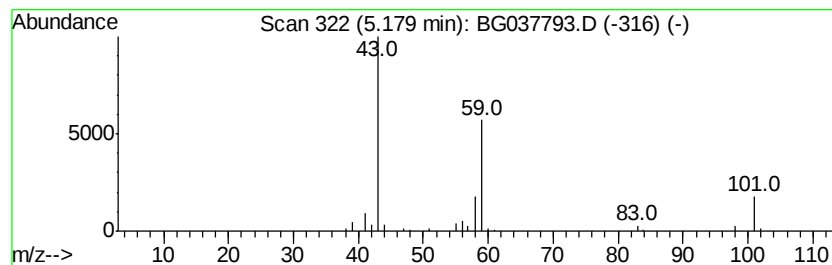
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	3.84 ng	67106	1,4-Dichlorobenzene-d4	8.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	45
3			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	33
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16



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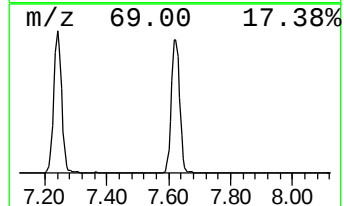
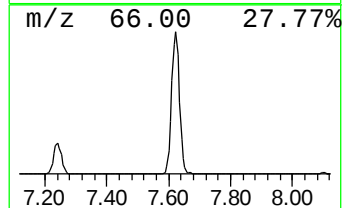
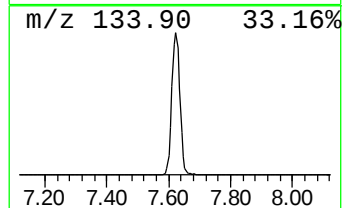
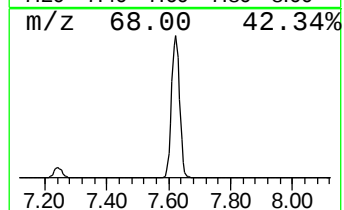
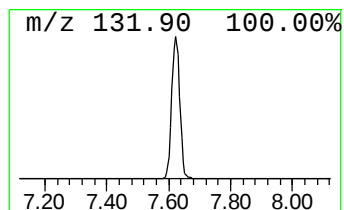
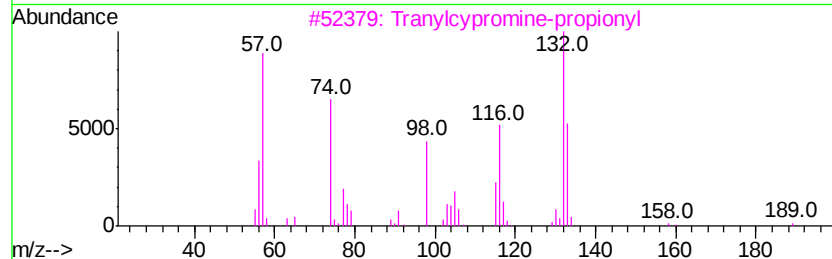
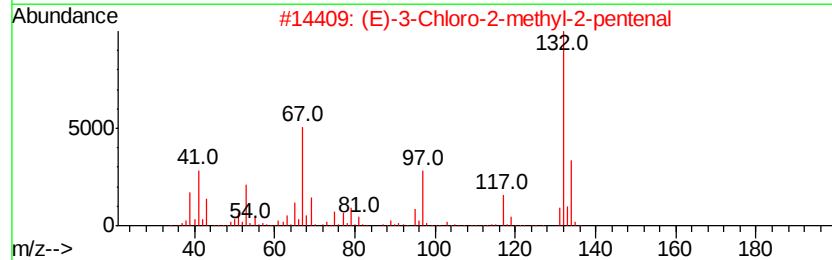
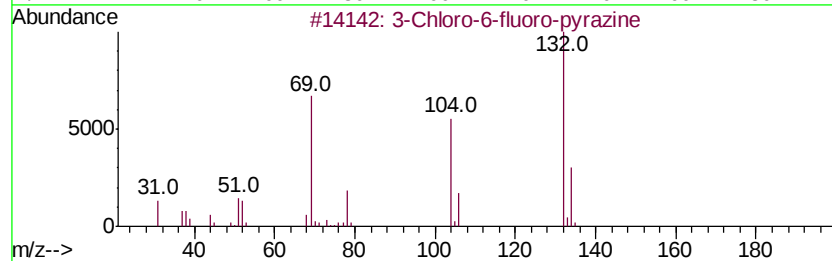
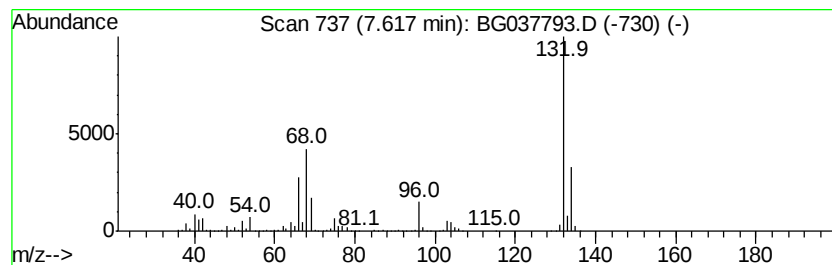
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	96.82 ng	1693110	1,4-Dichlorobenzene-d4	8.09

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		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	12
4		Amphetaminil	250	C17H18N2	017590-01-1	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	3.8	ng	67106	1	8.09	349759	20.0
unknown7.62	7.62	96.8	ng	1693110	1	8.09	349759	20.0