

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110719\
 Data File : BG043540.D
 Acq On : 7 Nov 2019 15:38
 Operator : JU
 Sample : PB124565BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB124565BL

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-BG102119.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	5.115	339	345	352	rBV	97834	156647	5.41%	1.258%
2	5.608	421	429	447	rBV	500819	880884	30.44%	7.072%
3	7.218	693	703	723	rBV	488879	943378	32.60%	7.573%
4	7.559	752	761	774	rBV	524110	957640	33.09%	7.688%
5	8.006	830	837	847	rBV	105458	186095	6.43%	1.494%
6	8.329	884	892	903	rBV	413445	757347	26.17%	6.080%
7	9.169	1027	1035	1050	rBV	272993	552106	19.08%	4.432%
8	10.814	1308	1315	1326	rBV	105914	220902	7.63%	1.773%
9	13.258	1724	1731	1746	rBV	842635	1413756	48.85%	11.349%
10	14.639	1957	1966	1978	rBV	203970	348500	12.04%	2.798%
11	16.131	2212	2220	2234	rBV	772169	1328399	45.90%	10.664%
12	17.383	2427	2433	2448	rBV	254738	468836	16.20%	3.764%
13	19.997	2871	2878	2894	rBV	2087616	2893839	100.00%	23.231%
14	21.654	3153	3160	3178	rBV2	315171	634800	21.94%	5.096%
15	23.117	3405	3409	3417	rVB6	14972	34678	1.20%	0.278%
16	23.916	3540	3545	3552	rVB7	15993	32005	1.11%	0.257%
17	24.845	3692	3703	3715	rBV2	205682	646819	22.35%	5.193%

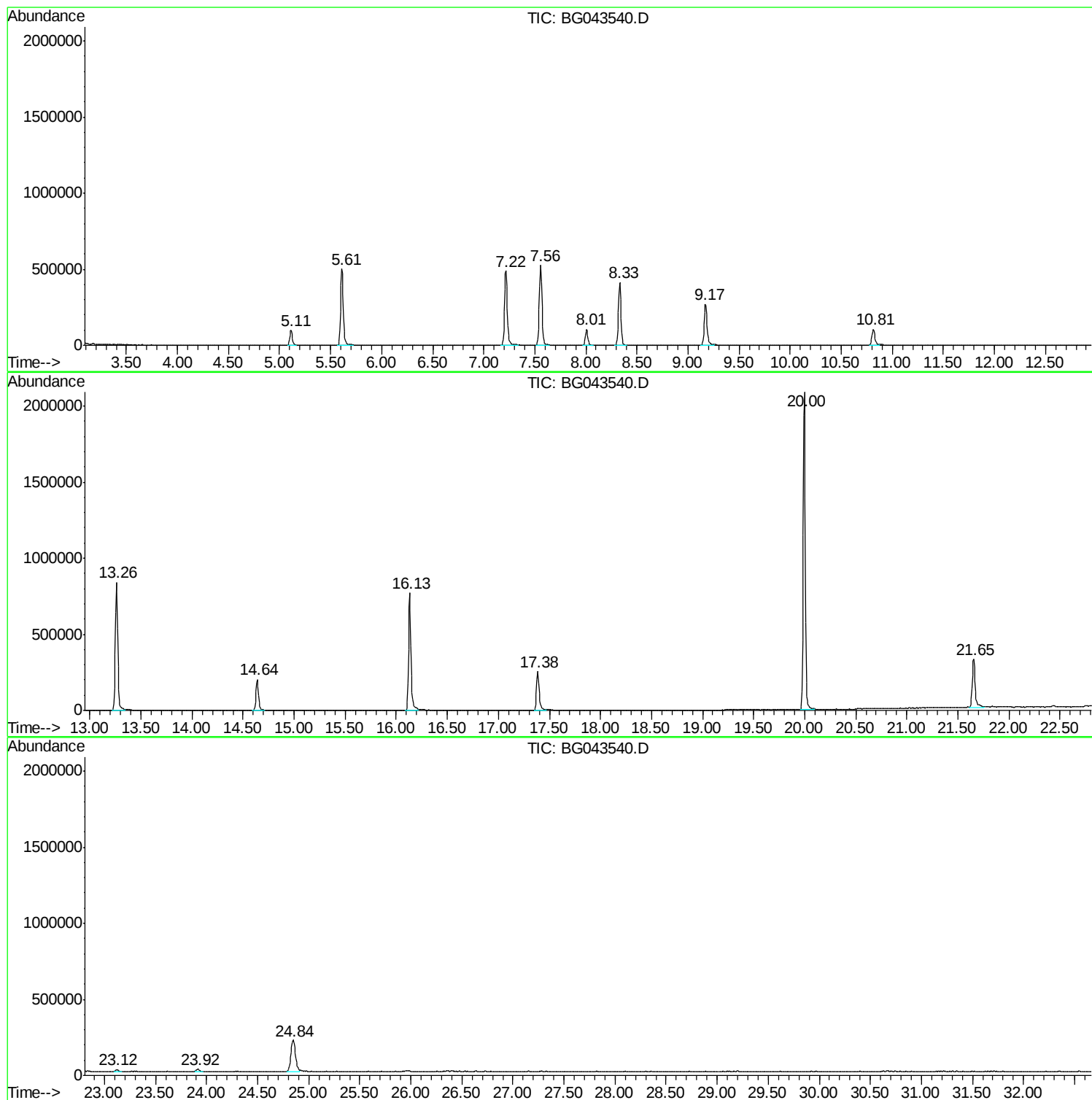
Sum of corrected areas: 12456631

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.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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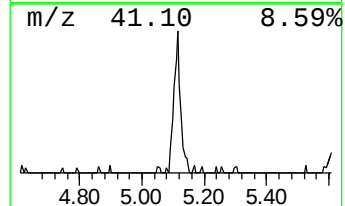
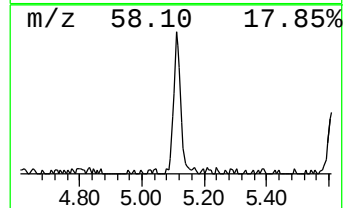
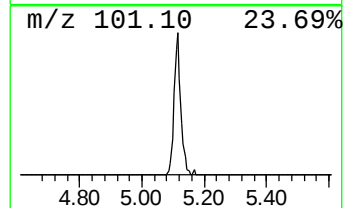
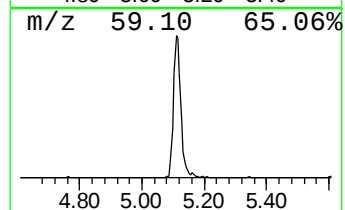
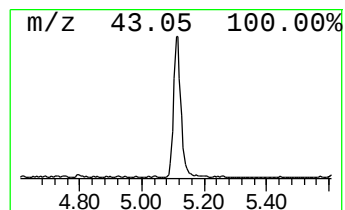
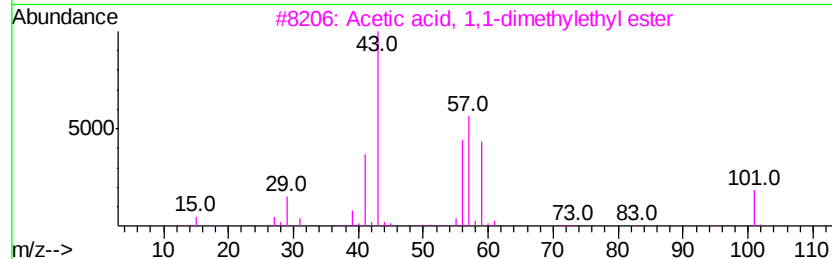
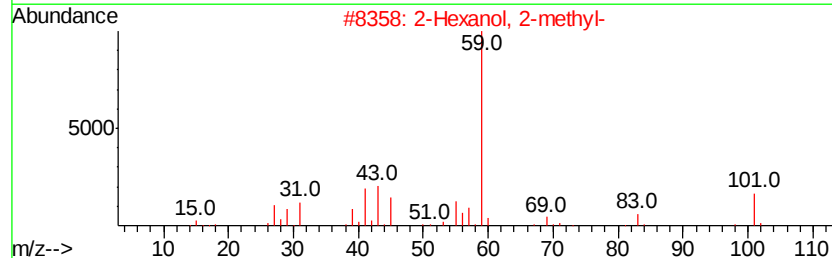
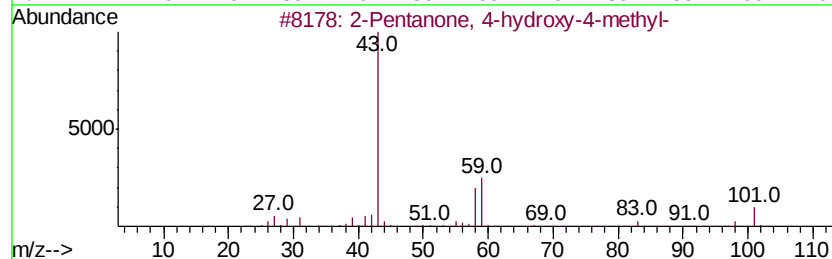
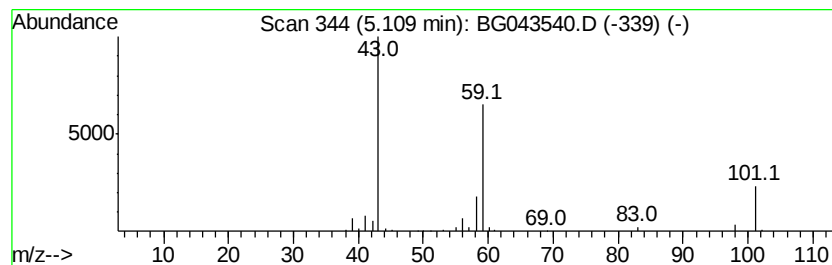
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	16.84 ng	156647	1,4-Dichlorobenzene-d4	8.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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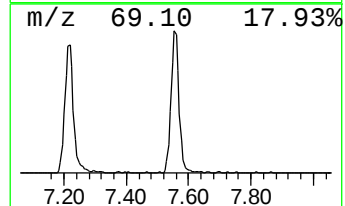
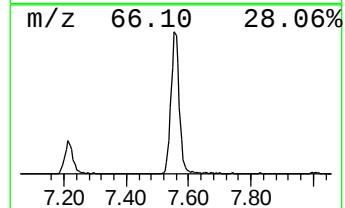
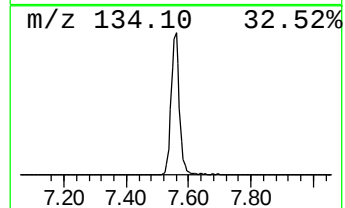
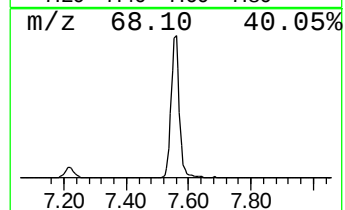
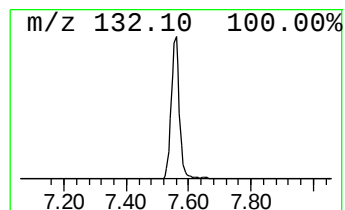
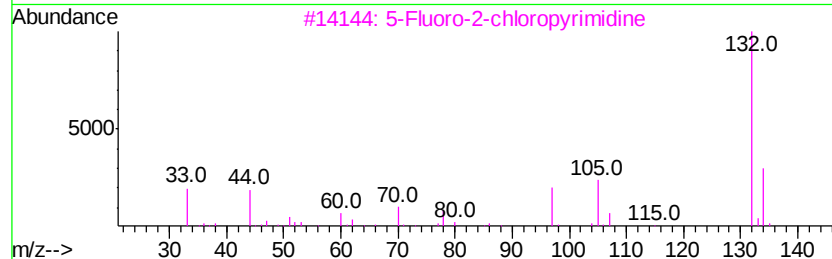
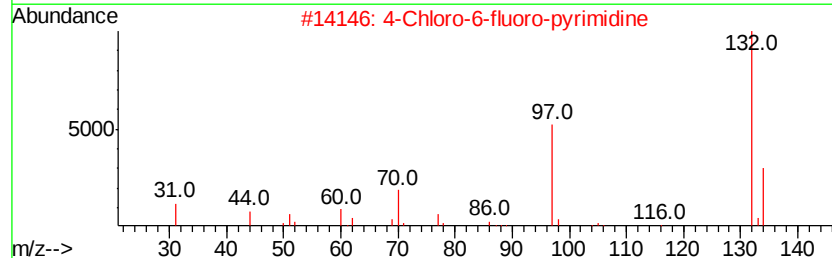
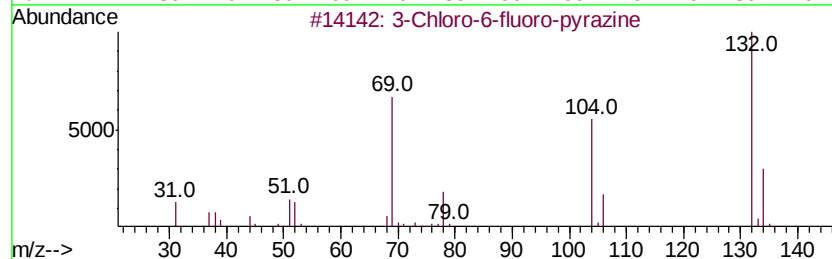
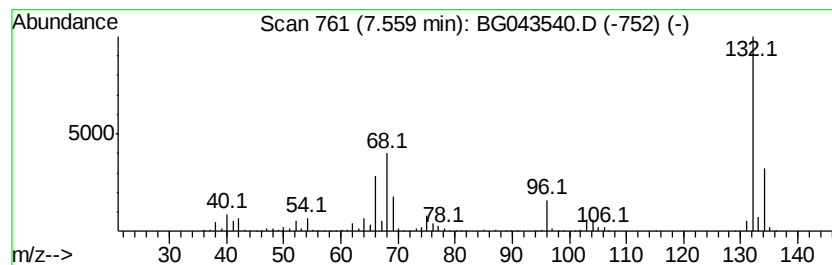
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Peak Number 2 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	102.92 ng	957640	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	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...	5.11	16.8	ng	156647	1	8.01	186095	20.0
unknown7.56	7.56	102.9	ng	957640	1	8.01	186095	20.0