

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040523\
 Data File : BG057005.D
 Acq On : 6 Apr 2023 00:48
 Operator : CG/JU
 Sample : 02126-02 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RS-B

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-BG033023.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG057005.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.441	318	324	332	rBV	25169	38984	11.55%	1.614%
2	5.922	399	406	415	rBB	60578	103092	30.54%	4.268%
3	7.544	675	682	692	rBB	67866	119176	35.31%	4.934%
4	8.413	824	830	837	rBB	76855	134726	39.92%	5.578%
5	9.588	1023	1030	1039	rVB2	40824	73903	21.90%	3.060%
6	11.256	1307	1314	1321	rBV	103397	186028	55.12%	7.702%
7	13.653	1715	1722	1729	rBV	117233	169546	50.23%	7.020%
8	14.763	1904	1911	1917	rBV2	2452	4173	1.24%	0.173%
9	15.028	1950	1956	1963	rVB	162868	253924	75.23%	10.513%
10	16.361	2177	2183	2189	rBV	3951	5971	1.77%	0.247%
11	16.508	2203	2208	2216	rVB	79383	122913	36.42%	5.089%
12	17.583	2388	2391	2396	rBV4	2623	4326	1.28%	0.179%
13	17.771	2416	2423	2427	rBV	184950	287471	85.17%	11.902%
14	17.812	2427	2430	2437	rVB2	26767	36767	10.89%	1.522%
15	17.900	2441	2445	2447	rBV3	6200	8986	2.66%	0.372%
16	18.153	2484	2488	2489	rBV4	3113	4217	1.25%	0.175%
17	18.482	2541	2544	2549	rBV6	2952	5482	1.62%	0.227%
18	18.605	2562	2565	2566	rBV3	6281	6234	1.85%	0.258%
19	18.635	2569	2570	2575	rVV5	7312	6325	1.87%	0.262%
20	18.676	2575	2577	2582	rVB5	8369	9302	2.76%	0.385%
21	18.828	2597	2603	2606	rBV3	11283	19753	5.85%	0.818%
22	19.069	2642	2644	2646	rVB3	5627	5157	1.53%	0.214%
23	19.116	2650	2652	2656	rVB5	9251	10107	2.99%	0.418%
24	19.146	2656	2657	2658	rBV	5319	3552	1.05%	0.147%
25	19.510	2717	2719	2720	rBV2	5864	3412	1.01%	0.141%
26	19.527	2720	2722	2723	rVV2	5417	4289	1.27%	0.178%
27	19.798	2763	2768	2772	rBV	86336	116854	34.62%	4.838%
28	20.156	2825	2829	2835	rVB	87346	128214	37.99%	5.309%
29	20.332	2855	2859	2864	rBV	163188	204817	60.68%	8.480%
30	22.089	3151	3158	3164	rBV2	154849	337526	100.00%	13.975%

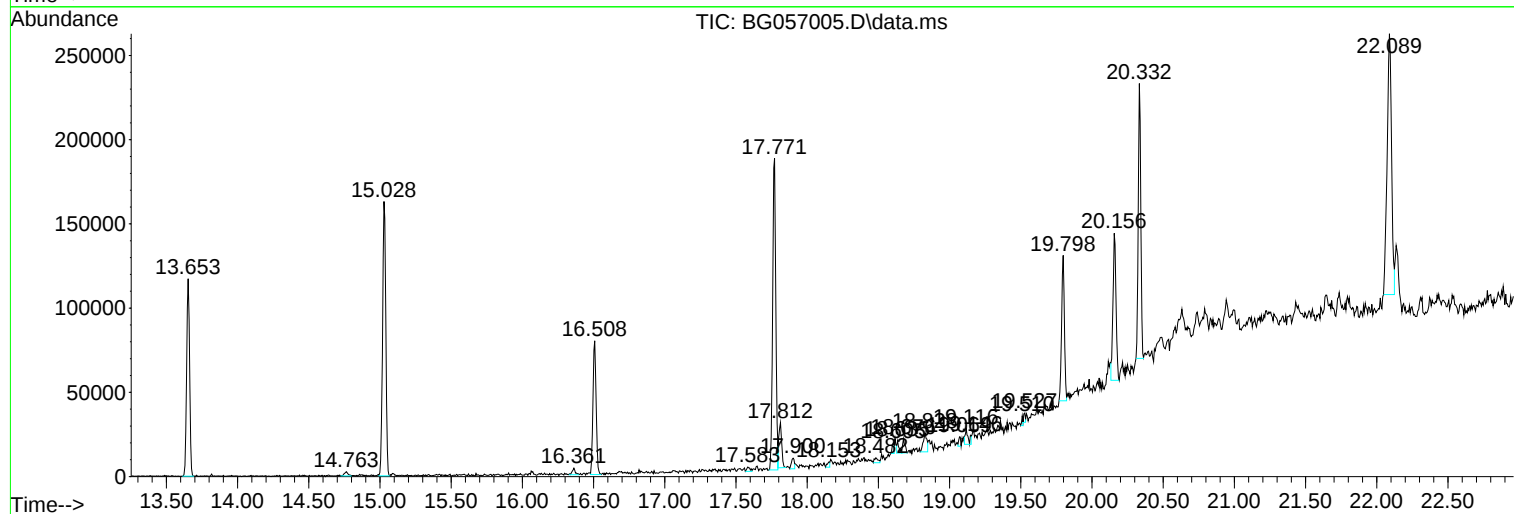
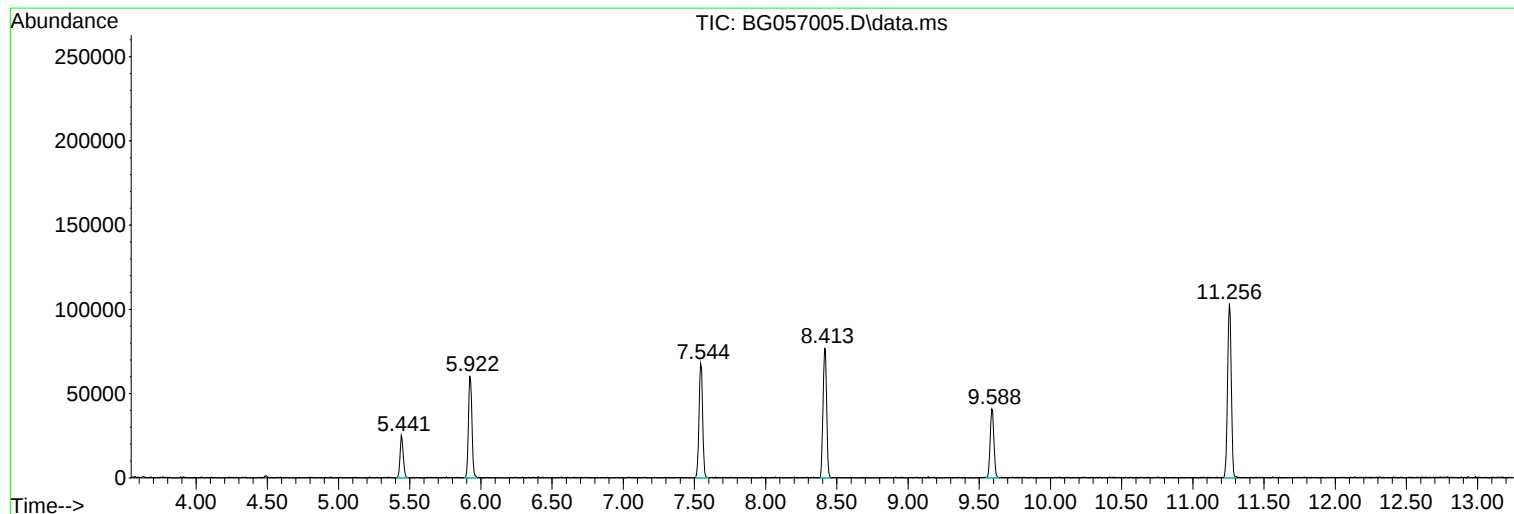
Sum of corrected areas: 2415227

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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.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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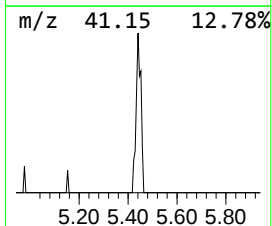
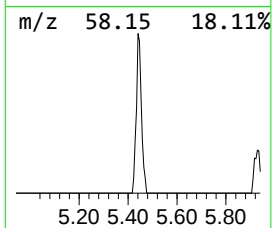
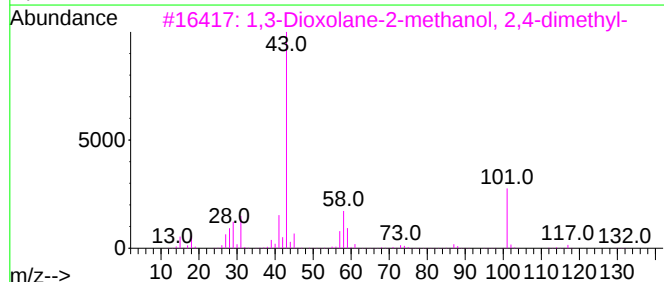
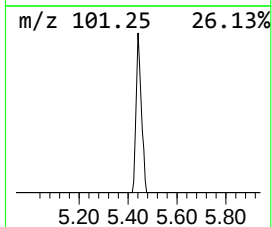
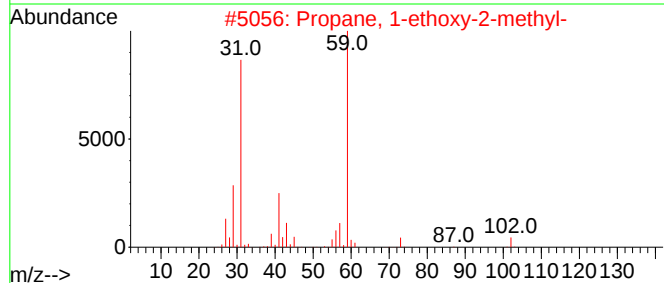
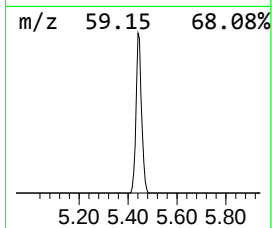
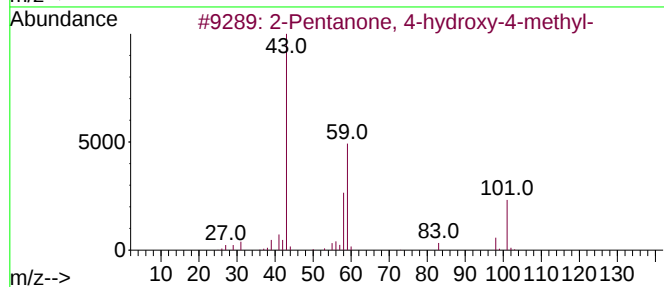
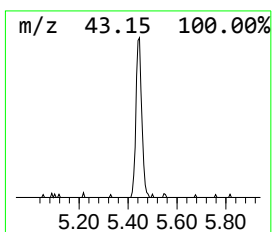
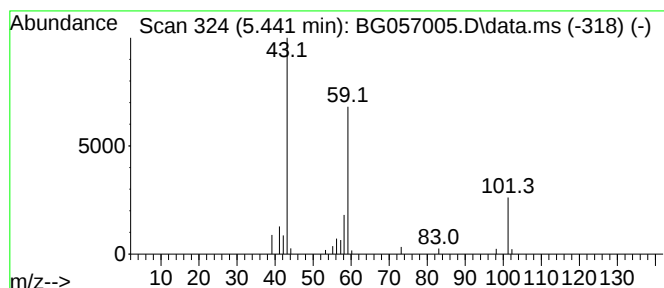
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.441	5.79 ng	38984	1,4-Dichlorobenzene-d4	8.413

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	38		
3	1,3-Dioxolane-2-methanol, 2,4-di...	132	C6H12O3	053951-43-2	28		
4	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	23		
5	N-(n-Propyl)acetamide	101	C5H11NO	005331-48-6	9		



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.441	5.8	ng	38984	1	8.413	134726	20.0