

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
 Data File : BN022717.D
 Acq On : 17 Nov 2022 13:24
 Operator : CG/JU
 Sample : N5622-11
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGCB3

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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_N\Methods\SFAM-EPA-BN111522.M
 Title : SVOA CALIBRATION

Signal : TIC: BN022717.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.181	29	33	44	rVB	12995	23197	1.00%	0.128%
2	3.628	106	109	126	rBV	54470	111702	4.79%	0.614%
3	3.852	143	147	152	rVB2	5322	8604	0.37%	0.047%
4	3.952	160	164	171	rVB3	4200	6994	0.30%	0.038%
5	4.923	324	329	340	rBV	35716	58474	2.51%	0.322%
6	7.028	681	687	698	rBV	74525	124491	5.34%	0.685%
7	7.193	707	715	726	rBV	225945	362110	15.54%	1.991%
8	7.387	741	748	757	rBV	221286	363011	15.58%	1.996%
9	7.858	822	828	837	rVB	199821	313118	13.44%	1.722%
10	8.587	944	952	965	rBV	197466	337126	14.47%	1.854%
11	9.034	1022	1028	1039	rBV	278570	470191	20.18%	2.585%
12	9.763	1142	1152	1162	rBV	231311	387447	16.63%	2.130%
13	10.305	1237	1244	1254	rBV	307520	523749	22.48%	2.880%
14	10.675	1300	1307	1314	rBV	298585	492132	21.12%	2.706%
15	10.828	1326	1333	1348	rBV	316028	552302	23.70%	3.037%
16	11.857	1501	1508	1512	rBV8	1703	3789	0.16%	0.021%
17	12.281	1576	1580	1586	rBV2	4173	6494	0.28%	0.036%
18	12.875	1676	1681	1683	rBV5	1715	2768	0.12%	0.015%
19	13.787	1829	1836	1842	rVB5	2203	4186	0.18%	0.023%
20	13.946	1856	1863	1872	rBV	730256	990806	42.52%	5.448%
21	14.222	1903	1910	1922	rBV	741140	1084508	46.55%	5.963%
22	14.528	1956	1962	1969	rBV	484978	705857	30.29%	3.881%
23	14.757	1995	2001	2017	rBV	68756	128207	5.50%	0.705%
24	14.963	2034	2036	2039	rVB4	2166	2339	0.10%	0.013%
25	15.528	2125	2132	2139	rBV	1007445	1416130	60.78%	7.787%
26	15.657	2148	2154	2165	rBV	371140	509855	21.88%	2.803%
27	15.769	2169	2173	2177	rVB4	3820	5414	0.23%	0.030%
28	17.287	2425	2431	2439	rBV	630602	871818	37.42%	4.794%
29	17.387	2442	2448	2458	rBV	1144287	1616608	69.38%	8.889%
30	17.492	2463	2466	2469	rVB4	2115	2383	0.10%	0.013%
31	18.181	2578	2583	2589	rBV2	43258	67826	2.91%	0.373%
32	18.263	2595	2597	2600	rVB3	2540	2727	0.12%	0.015%
33	19.251	2761	2765	2773	rBV2	12269	18952	0.81%	0.104%
34	19.328	2773	2778	2783	rBV	13665	22495	0.97%	0.124%
35	19.463	2797	2801	2812	rBV5	31394	57527	2.47%	0.316%
36	19.616	2823	2827	2829	rBV4	4511	6446	0.28%	0.035%

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Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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

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37	19.692	2834	2840	2851	rVV	1441346	1901909	81.63%	10.458%
38	21.492	3141	3146	3154	rBV	839479	1069973	45.92%	5.883%
39	21.616	3165	3167	3175	rVB4	21827	37049	1.59%	0.204%
40	23.745	3522	3529	3542	rVB2	1094939	2329956	100.00%	12.811%
41	23.904	3549	3556	3567	rVB2	577583	1185865	50.90%	6.521%

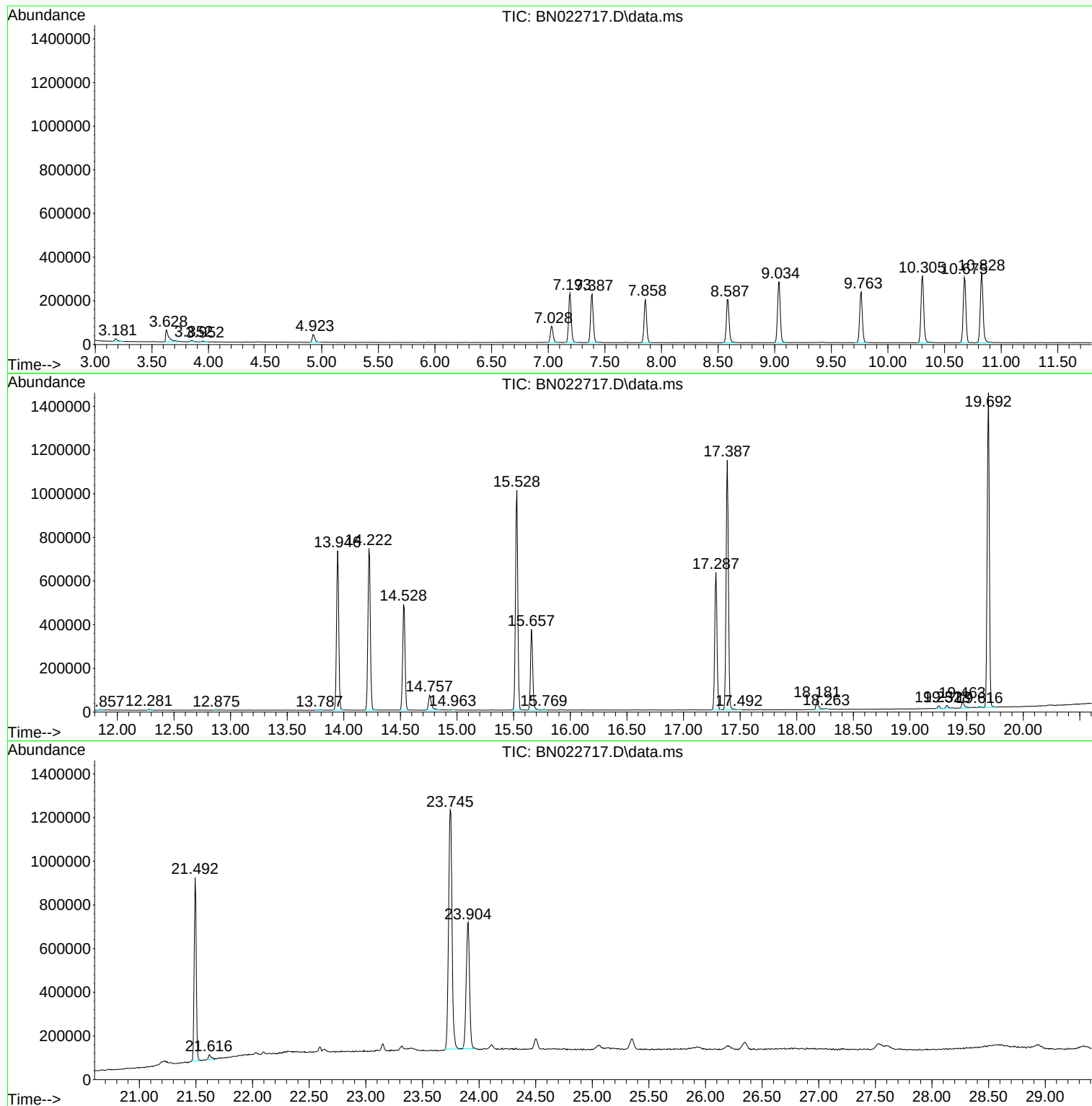
Sum of corrected areas: 18186535

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
Quant Title : SVOA CALIBRATION

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



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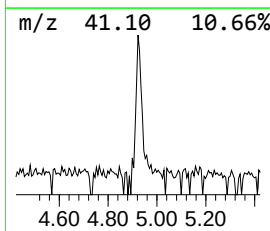
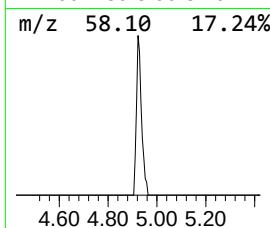
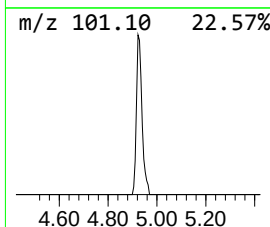
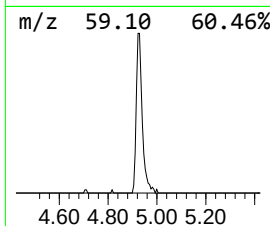
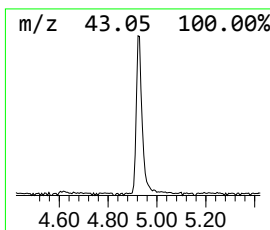
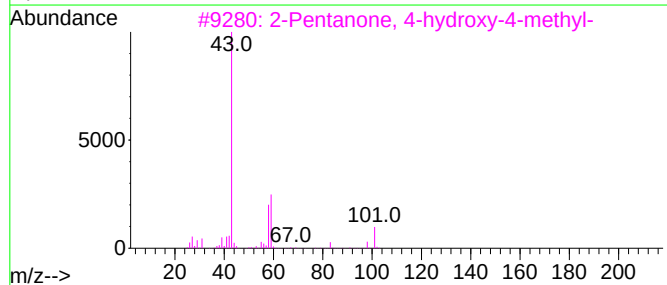
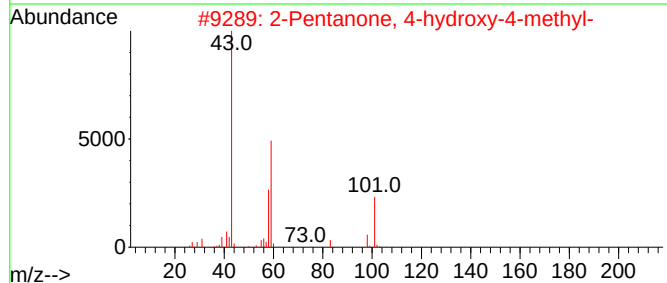
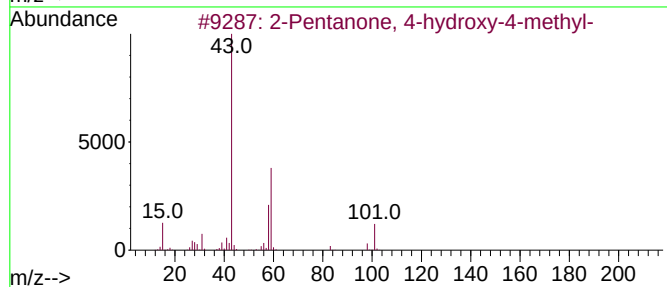
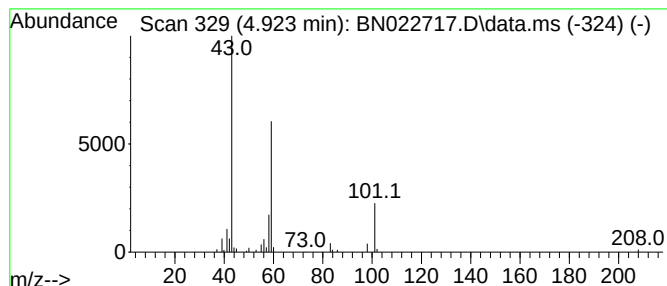
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.923	3.73 ng/ul	58474	1,4-Dichlorobenzene-d4	7.858

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38		
5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.923	3.7	ng/u1	58474	1	7.858	313118	20.0