

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111522\
 Data File : BN022688.D
 Acq On : 16 Nov 2022 15:43
 Operator : CG/JU
 Sample : N5622-06
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGCA8

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: BN022688.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.193	31	35	40	rBV	17427	23505	1.08%	0.118%
2	3.634	107	110	130	rBV	60460	116599	5.37%	0.585%
3	3.863	142	149	156	rBV	8763	16956	0.78%	0.085%
4	3.957	159	165	177	rVB	17030	27669	1.27%	0.139%
5	4.934	326	331	343	rBV	39747	60204	2.77%	0.302%
6	7.045	684	690	700	rVB	87471	139813	6.44%	0.702%
7	7.204	710	717	729	rBV	275707	432801	19.94%	2.172%
8	7.398	743	750	759	rBV	280805	450908	20.78%	2.263%
9	7.875	824	831	840	rVB	258089	418148	19.27%	2.099%
10	8.604	949	955	966	rBV	243728	386008	17.79%	1.937%
11	9.057	1023	1032	1043	rBV	348908	574384	26.47%	2.883%
12	9.175	1050	1052	1056	rVB4	2428	2886	0.13%	0.014%
13	9.781	1148	1155	1167	rBV	293183	475493	21.91%	2.386%
14	10.328	1241	1248	1261	rBV	395189	663927	30.59%	3.332%
15	10.698	1304	1311	1320	rBV	370299	625281	28.81%	3.138%
16	10.851	1329	1337	1353	rBV	376525	678642	31.27%	3.406%
17	12.310	1579	1585	1590	rVB3	4735	8537	0.39%	0.043%
18	13.974	1861	1868	1882	rVB	877572	1186904	54.69%	5.957%
19	14.251	1908	1915	1929	rBV	888144	1299728	59.89%	6.523%
20	14.557	1961	1967	1977	rBV	594569	858496	39.56%	4.309%
21	14.792	2002	2007	2025	rBV	59338	127846	5.89%	0.642%
22	14.992	2038	2041	2046	rVB5	4773	5997	0.28%	0.030%
23	15.557	2130	2137	2146	rBV	1191558	1678335	77.33%	8.423%
24	15.686	2153	2159	2174	rVB	401868	554595	25.55%	2.783%
25	15.798	2174	2178	2181	rVB2	4018	4671	0.22%	0.023%
26	16.763	2337	2342	2347	rBV7	1307	2449	0.11%	0.012%
27	17.315	2430	2436	2446	rBV	687402	997358	45.95%	5.005%
28	17.421	2447	2454	2466	rVV	1409700	2042825	94.12%	10.252%
29	17.527	2470	2472	2475	rVB3	4103	4228	0.19%	0.021%
30	18.215	2585	2589	2594	rVB2	8473	10278	0.47%	0.052%
31	19.286	2767	2771	2779	rBV3	14662	21550	0.99%	0.108%
32	19.362	2779	2784	2788	rBV	13788	19286	0.89%	0.097%
33	19.504	2804	2808	2811	rBV4	6950	8865	0.41%	0.044%
34	19.727	2840	2846	2855	rBV	1644585	2170353	100.00%	10.892%
35	21.527	3147	3152	3161	rBV	765207	992203	45.72%	4.980%
36	23.803	3532	3539	3553	rVB2	939304	1978263	91.15%	9.928%

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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	23.962	3560	3566	3575	rVB2	429777	859537	39.60%	4.314%
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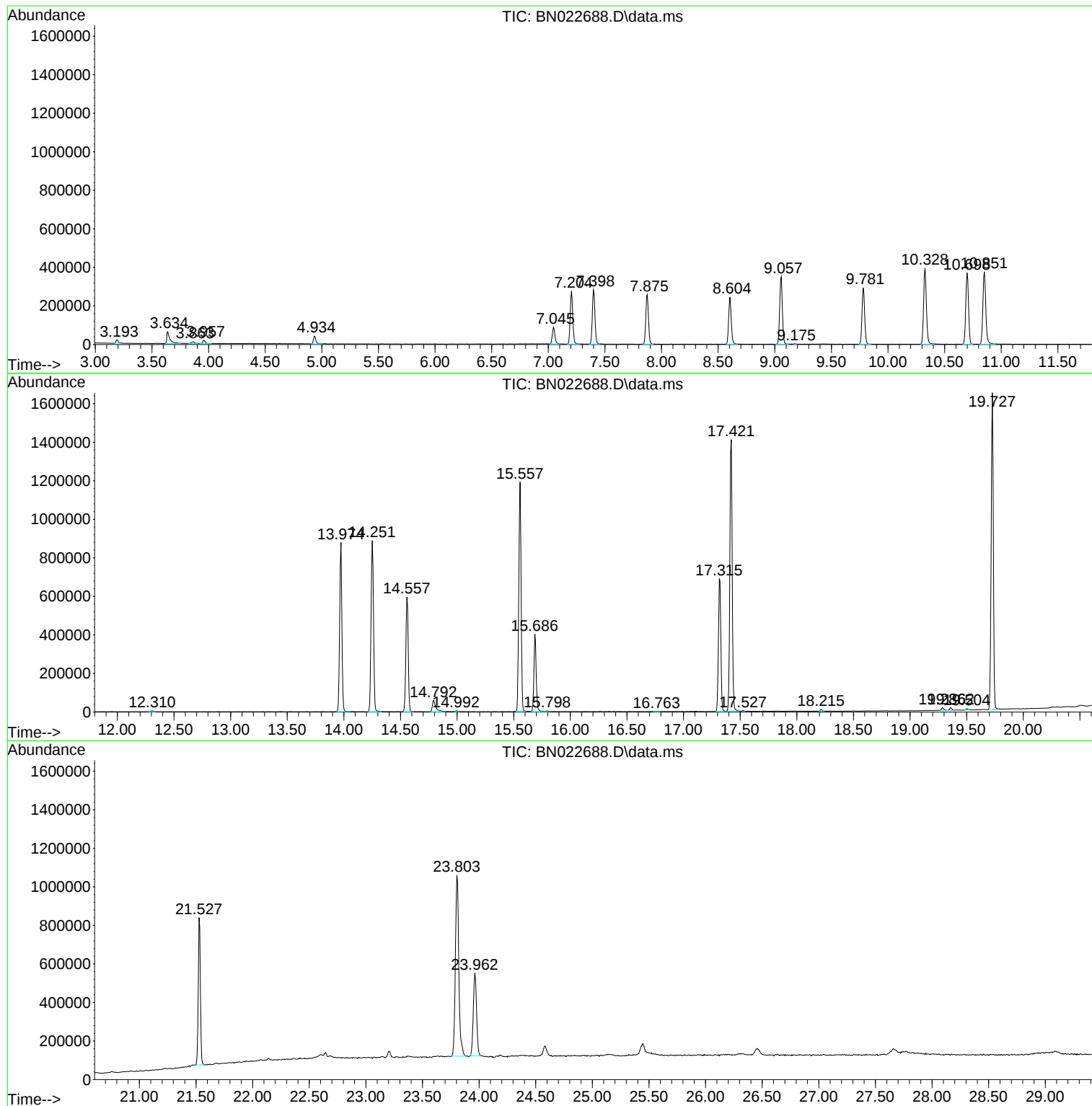
Sum of corrected areas: 19925528

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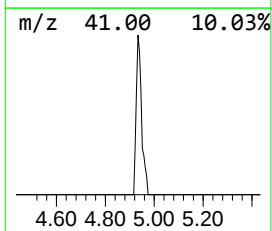
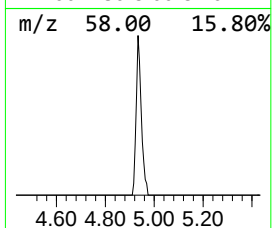
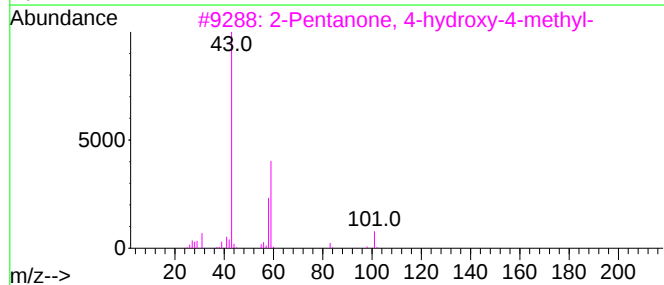
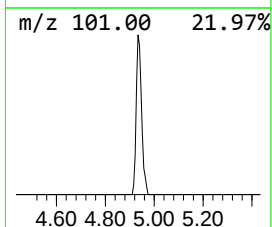
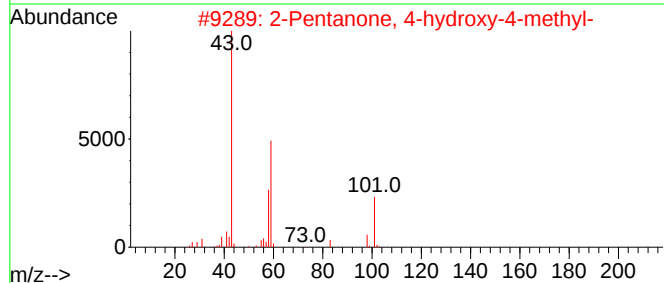
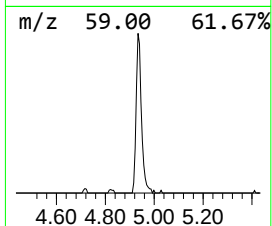
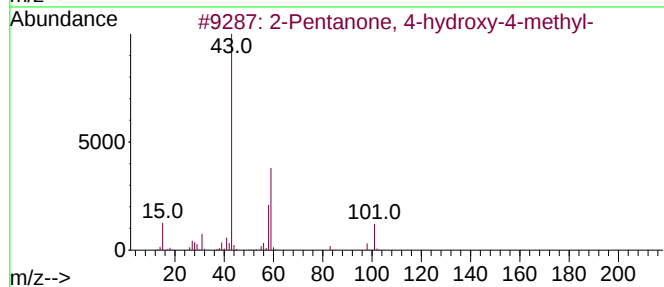
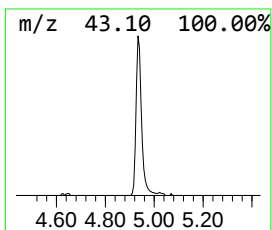
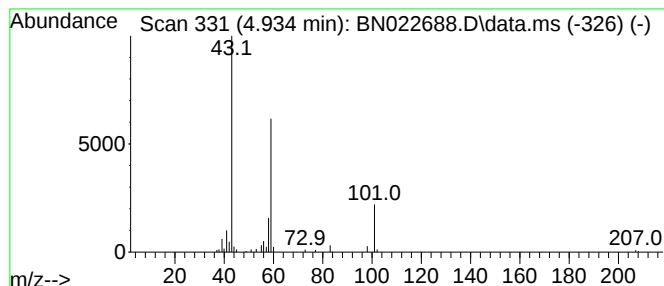
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.934	2.88 ng/ul	60204	1,4-Dichlorobenzene-d4	7.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	39		
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	36		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.934	2.9	ng/u1	60204	1	7.875	418148	20.0