

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111522\
 Data File : BN022686.D
 Acq On : 16 Nov 2022 14:29
 Operator : CG/JU
 Sample : N5622-04
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGCA6

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: BN022686.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	44	rBV	15978	23601	1.11%	0.112%
2	3.516	87	90	95	rBV3	1724	2667	0.13%	0.013%
3	3.634	107	110	132	rBV	64770	122703	5.77%	0.582%
4	3.858	143	148	157	rVB	10115	19368	0.91%	0.092%
5	3.958	162	165	172	rVB	5592	7902	0.37%	0.037%
6	4.934	326	331	339	rVV	49142	73376	3.45%	0.348%
7	7.046	684	690	700	rVB	95114	150193	7.06%	0.712%
8	7.205	710	717	729	rBV	332189	523277	24.61%	2.482%
9	7.399	743	750	759	rBV	330673	529994	24.92%	2.514%
10	7.475	761	763	769	rVB4	1812	2262	0.11%	0.011%
11	7.875	824	831	842	rBV	273180	443190	20.84%	2.102%
12	8.604	948	955	965	rBV	267301	436333	20.52%	2.069%
13	9.057	1025	1032	1043	rBV	419207	693952	32.63%	3.291%
14	9.446	1091	1098	1102	rVB4	1335	2866	0.13%	0.014%
15	9.781	1149	1155	1170	rVB	357873	571157	26.86%	2.709%
16	10.328	1241	1248	1261	rBV	461449	783210	36.83%	3.715%
17	10.698	1304	1311	1319	rBV	394672	663297	31.19%	3.146%
18	10.851	1328	1337	1357	rBV	407197	724910	34.09%	3.438%
19	12.310	1579	1585	1590	rBV	5860	9694	0.46%	0.046%
20	13.445	1774	1778	1783	rVB4	1610	2883	0.14%	0.014%
21	13.975	1861	1868	1880	rBV	908958	1284825	60.41%	6.094%
22	14.251	1908	1915	1930	rBV	1023605	1491857	70.15%	7.076%
23	14.557	1960	1967	1976	rBV	633927	894565	42.06%	4.243%
24	14.792	2002	2007	2026	rBV	50784	114005	5.36%	0.541%
25	14.992	2038	2041	2047	rVB3	7046	9559	0.45%	0.045%
26	15.557	2130	2137	2151	rBV	1288701	1841080	86.57%	8.732%
27	15.686	2153	2159	2173	rVV	396556	529725	24.91%	2.512%
28	15.792	2173	2177	2181	rVB	4095	5582	0.26%	0.026%
29	17.316	2431	2436	2447	rBV	761109	1047442	49.25%	4.968%
30	17.416	2447	2453	2469	rVV2	1321220	1987895	93.47%	9.428%
31	18.210	2585	2588	2595	rVB2	10258	13642	0.64%	0.065%
32	19.286	2767	2771	2777	rBV2	14819	19203	0.90%	0.091%
33	19.357	2780	2783	2788	rBV2	12204	15773	0.74%	0.075%
34	19.498	2804	2807	2812	rBV4	9477	13352	0.63%	0.063%
35	19.727	2839	2846	2855	rBV	1570462	2126676	100.00%	10.086%
36	21.527	3147	3152	3161	rBV	831368	1034008	48.62%	4.904%

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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	23.804	3531	3539	3553	rVB2	929569	1939865	91.22%	9.200%
38	23.957	3559	3565	3578	rVB2	464448	928588	43.66%	4.404%

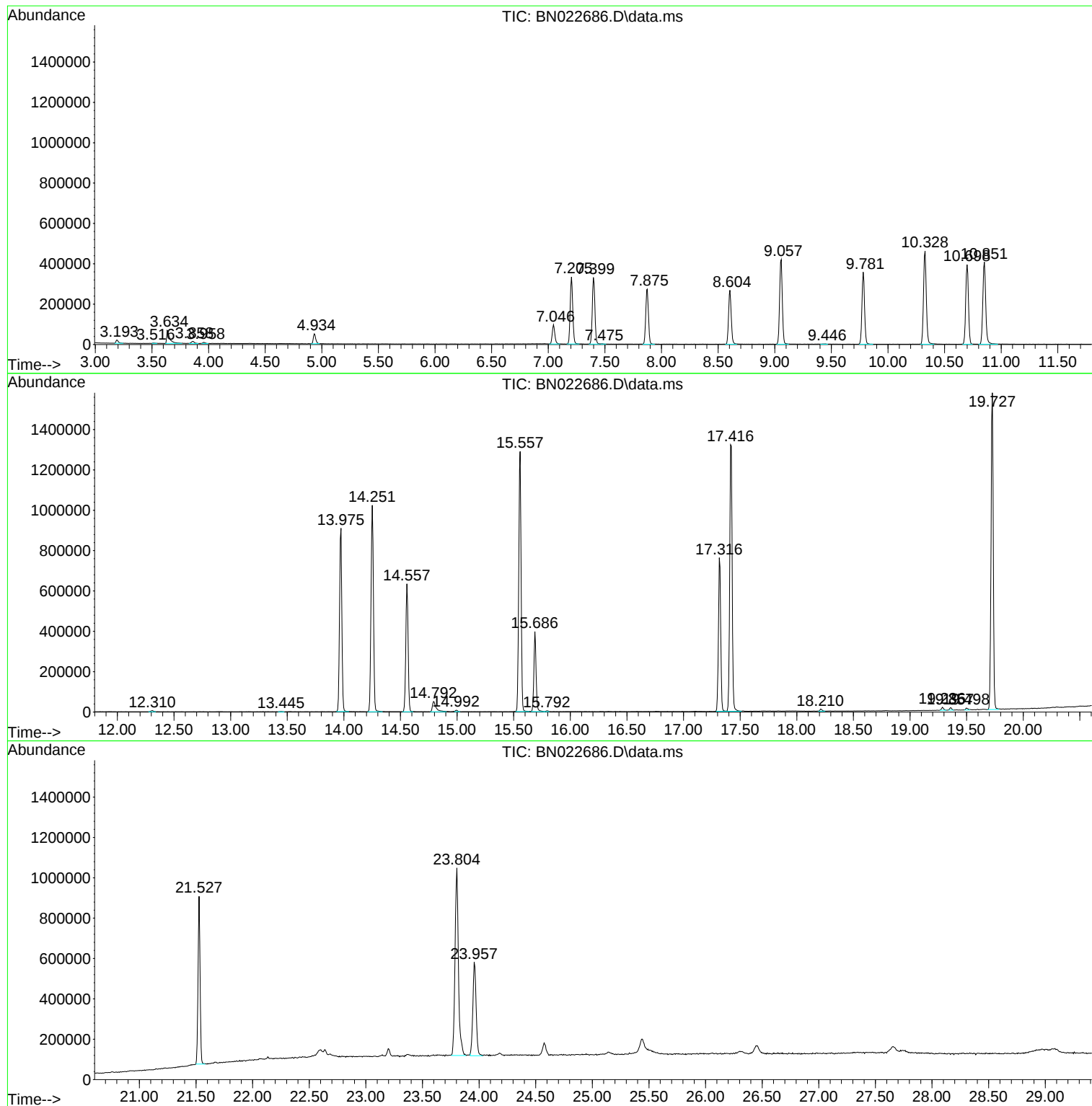
Sum of corrected areas: 21084477

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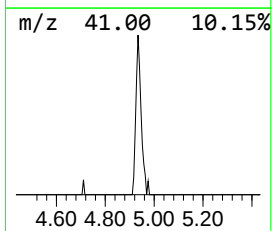
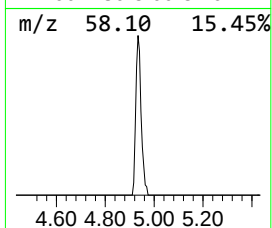
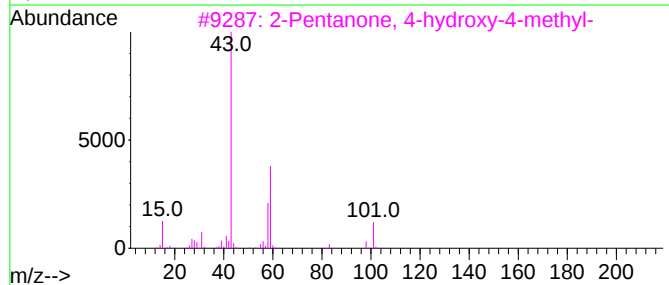
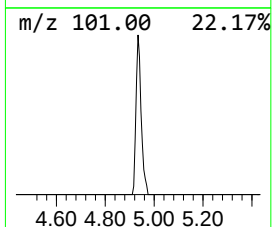
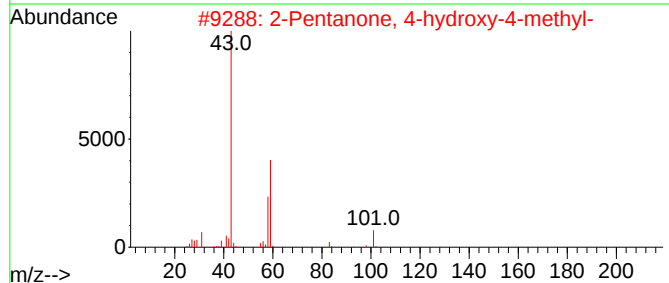
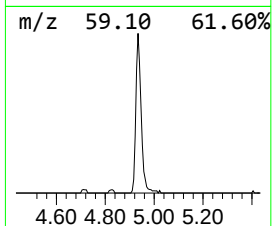
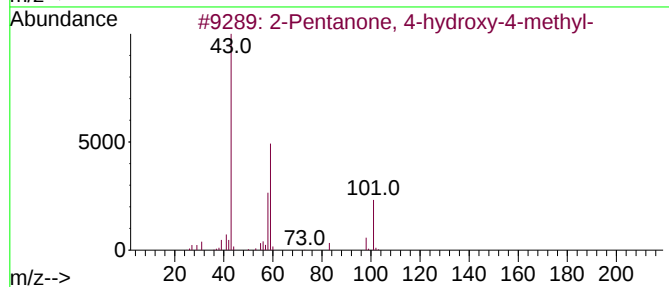
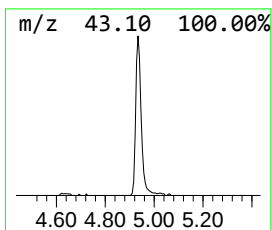
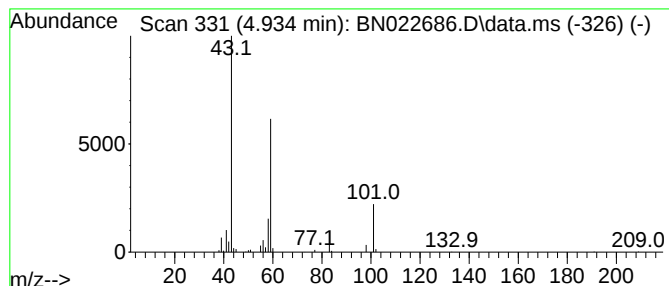
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.934	3.31 ng/ul	73376	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	33		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25		
5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	23		



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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-...	4.934	3.3	ng/u1	73376	1	7.875	443190	20.0