

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
 Data File : BN022727.D
 Acq On : 17 Nov 2022 19:33
 Operator : CG/JU
 Sample : N5622-16
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGCA1

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: BN022727.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	42	rVB	13370	19581	0.90%	0.106%
2	3.622	106	108	126	rBV	68280	130311	5.98%	0.704%
3	3.852	141	147	152	rVB2	6415	10232	0.47%	0.055%
4	3.946	160	163	167	rBV	3224	4213	0.19%	0.023%
5	4.922	325	329	339	rBV	28950	43785	2.01%	0.236%
6	7.028	681	687	702	rVB	61164	100169	4.60%	0.541%
7	7.187	707	714	728	rBV	213528	346445	15.91%	1.870%
8	7.381	741	747	758	rBV	202081	328701	15.10%	1.775%
9	7.857	820	828	839	rBV	210657	342578	15.73%	1.850%
10	8.581	945	951	965	rBV	177423	291646	13.39%	1.575%
11	9.034	1020	1028	1041	rBV	278977	458574	21.06%	2.476%
12	9.757	1145	1151	1164	rVB	216683	363750	16.70%	1.964%
13	10.298	1237	1243	1255	rBV	279557	484607	22.26%	2.616%
14	10.675	1300	1307	1315	rBV	340729	559784	25.71%	3.022%
15	10.828	1326	1333	1354	rBV	310117	558294	25.64%	3.014%
16	11.845	1502	1506	1512	rBV4	935	2387	0.11%	0.013%
17	11.916	1512	1518	1523	rVB2	1275	2785	0.13%	0.015%
18	12.287	1575	1581	1588	rVB2	3407	6652	0.31%	0.036%
19	13.945	1854	1863	1876	rVB	703163	1006591	46.23%	5.435%
20	14.222	1903	1910	1922	rBV2	761442	1083120	49.74%	5.848%
21	14.528	1955	1962	1972	rBV	586757	834219	38.31%	4.504%
22	14.757	1995	2001	2017	rBV	51736	97935	4.50%	0.529%
23	14.963	2031	2036	2040	rVB4	4559	5286	0.24%	0.029%
24	15.522	2125	2131	2140	rVB	1045977	1455693	66.85%	7.859%
25	15.657	2148	2154	2167	rVB	341252	480978	22.09%	2.597%
26	15.763	2169	2172	2176	rVB	2786	3393	0.16%	0.018%
27	17.280	2424	2430	2439	rBV	730806	1037959	47.67%	5.604%
28	17.380	2441	2447	2459	rVV	1206576	1696774	77.92%	9.161%
29	17.486	2462	2465	2471	rVB3	9414	11895	0.55%	0.064%
30	17.951	2541	2544	2548	rBV3	2180	3080	0.14%	0.017%
31	18.180	2578	2583	2587	rBV	6153	8104	0.37%	0.044%
32	19.251	2760	2765	2771	rBV4	12377	18420	0.85%	0.099%
33	19.321	2771	2777	2781	rBV2	11724	16389	0.75%	0.088%
34	19.463	2798	2801	2807	rBV7	5157	8036	0.37%	0.043%
35	19.610	2823	2826	2828	rBV3	3415	3411	0.16%	0.018%
36	19.686	2833	2839	2849	rBV2	1457679	1935374	88.88%	10.449%

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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	21.486	3140	3145	3155	rBV2	956036	1187538	54.54%	6.412%
38	23.133	3423	3425	3432	rVB3	40252	59208	2.72%	0.320%
39	23.733	3520	3527	3540	rVB2	1019314	2177495	100.00%	11.756%
40	23.892	3546	3554	3563	rVB2	592643	1224427	56.23%	6.611%
41	24.486	3652	3655	3663	rVB	62967	112031	5.14%	0.605%

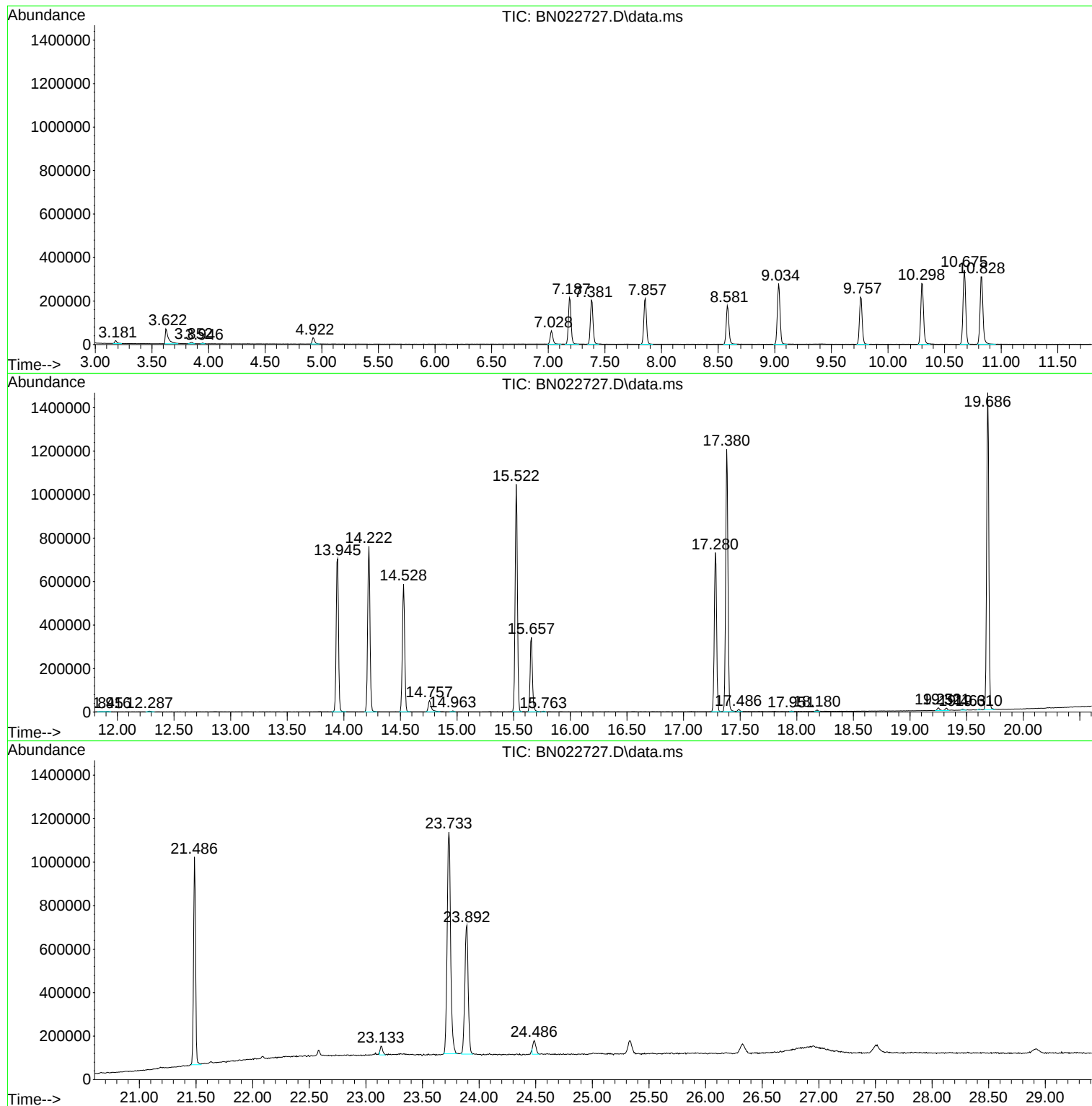
Sum of corrected areas: 18521850

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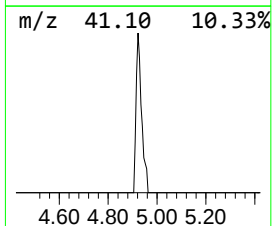
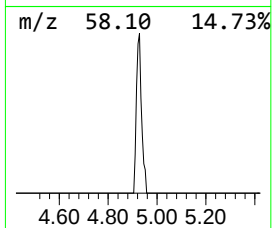
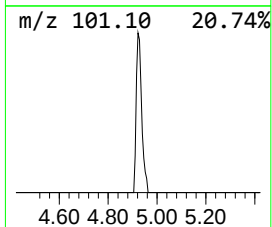
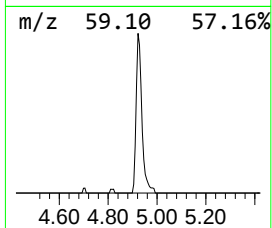
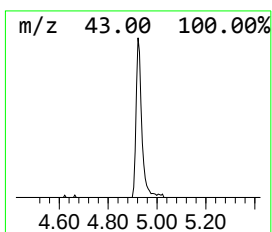
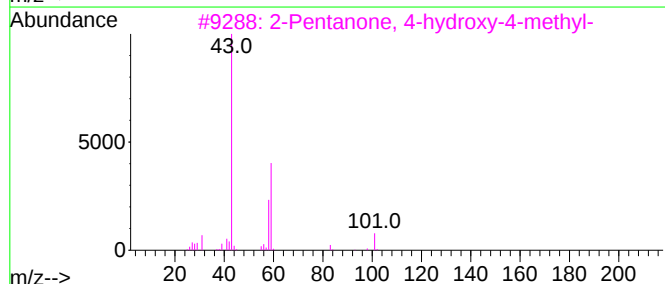
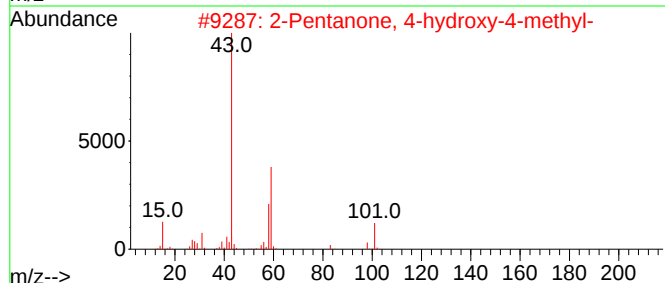
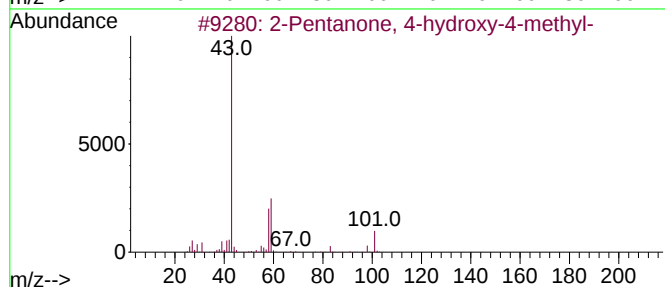
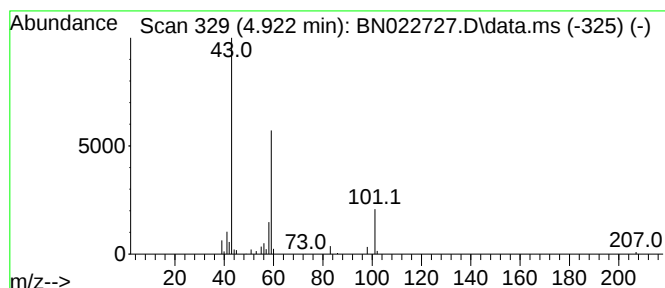
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
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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.
4.922	2.56 ng/ul	43785	1,4-Dichlorobenzene-d4	7.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42		
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38		
5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	32		



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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-...	4.922	2.6	ng/u1	43785	1	7.857	342578	20.0