

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101422\
 Data File : BP012141.D
 Acq On : 14 Oct 2022 17:00
 Operator : CG/JU
 Sample : N4982-15
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C1BC1

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_P\Methods\SFAM-EPA-BP101122.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012141.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.275	46	49	53	rVB	46859	52438	0.79%	0.093%
2	3.699	119	121	137	rBV	191490	304826	4.58%	0.543%
3	4.017	172	175	182	rVB	43429	57471	0.86%	0.102%
4	4.952	329	334	348	rVB	271420	410264	6.17%	0.731%
5	6.963	671	676	680	rBV	33472	49362	0.74%	0.088%
6	7.016	680	685	698	rVB	250003	437956	6.58%	0.780%
7	7.181	706	713	734	rVB	880297	1455507	21.88%	2.592%
8	7.375	739	746	757	rBV	880415	1413026	21.25%	2.516%
9	7.846	819	826	835	rBV	930090	1506035	22.64%	2.682%
10	7.910	835	837	844	rVB6	7712	13774	0.21%	0.025%
11	8.263	892	897	904	rBV3	6799	13102	0.20%	0.023%
12	8.563	941	948	965	rBV	645536	1117606	16.80%	1.990%
13	9.010	1017	1024	1038	rBV	945852	1623461	24.41%	2.891%
14	9.410	1086	1092	1098	rBV	51583	81072	1.22%	0.144%
15	9.734	1140	1147	1164	rBV	724941	1212251	18.23%	2.159%
16	10.275	1231	1239	1254	rBV	1035330	1816934	27.32%	3.236%
17	10.581	1287	1291	1295	rBV	4731	7008	0.11%	0.012%
18	10.651	1295	1303	1319	rVB	1310842	2215036	33.30%	3.945%
19	10.798	1319	1328	1346	rBV	535139	1045124	15.71%	1.861%
20	11.904	1511	1516	1523	rVB2	22763	35625	0.54%	0.063%
21	12.069	1539	1544	1550	rBV7	6261	12960	0.19%	0.023%
22	12.187	1554	1564	1570	rVV	64057	126420	1.90%	0.225%
23	12.245	1570	1574	1586	rVB	17568	40113	0.60%	0.071%
24	13.310	1748	1755	1760	rBV2	14994	25111	0.38%	0.045%
25	13.463	1776	1781	1786	rVB6	3902	7907	0.12%	0.014%
26	13.904	1849	1856	1870	rBV	1712835	2515308	37.82%	4.479%
27	14.187	1897	1904	1917	rBV	2241554	3302602	49.66%	5.881%
28	14.387	1933	1938	1942	rVB3	5530	8015	0.12%	0.014%
29	14.492	1949	1956	1970	rBV	1680102	2537294	38.15%	4.519%
30	14.716	1988	1994	2007	rBV	82585	210074	3.16%	0.374%
31	14.916	2024	2028	2039	rVB	8377	24322	0.37%	0.043%
32	15.492	2117	2126	2141	rBV	2757552	4074313	61.26%	7.256%
33	15.616	2141	2147	2162	rVV	668410	958426	14.41%	1.707%
34	15.728	2162	2166	2175	rVB3	14389	27739	0.42%	0.049%
35	17.251	2419	2425	2436	rBV	1881770	2776808	41.75%	4.945%
36	17.351	2436	2442	2458	rVV	2930074	4366019	65.64%	7.775%

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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

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37	18.139	2572	2576	2583	rBV	112804	171597	2.58%	0.306%
38	19.169	2748	2751	2756	rVB	34796	44709	0.67%	0.080%
39	19.233	2759	2762	2771	rBV	50085	87123	1.31%	0.155%
40	19.375	2782	2786	2792	rBV	101596	151307	2.27%	0.269%
41	19.592	2817	2823	2836	rBV	4375724	5721679	86.03%	10.190%
42	21.345	3116	3121	3129	rBV	2754359	3568481	53.65%	6.355%
43	23.610	3499	3506	3518	rVB2	3339130	6651078	100.00%	11.845%
44	23.763	3526	3532	3544	rVB2	1935616	3875230	58.26%	6.901%

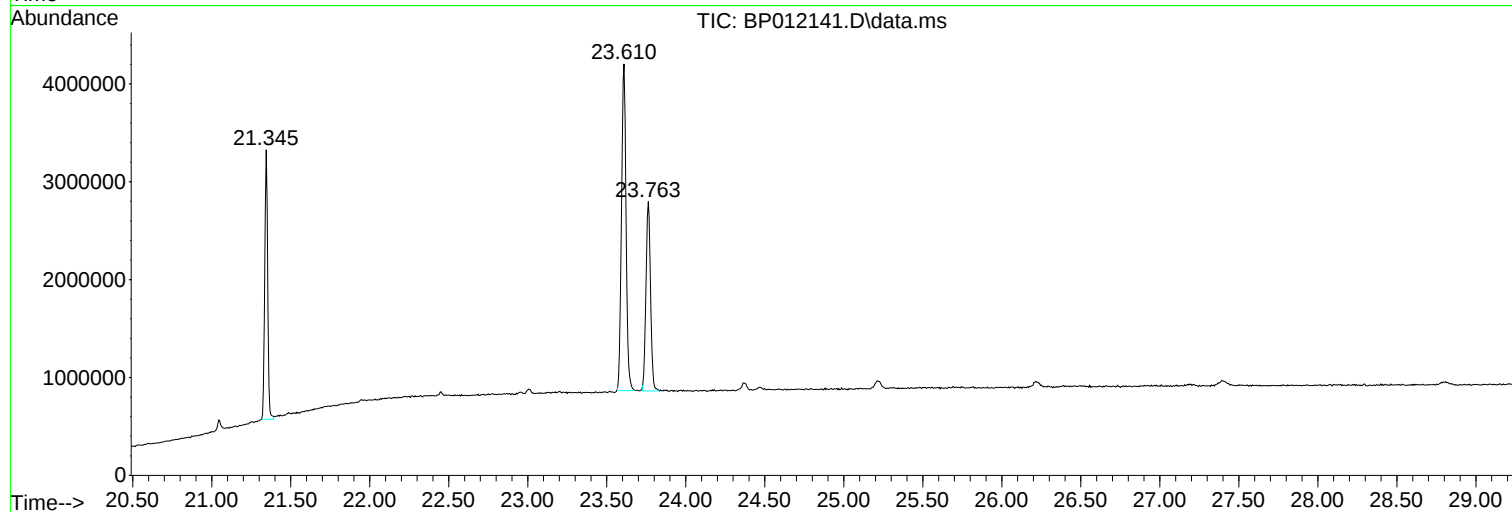
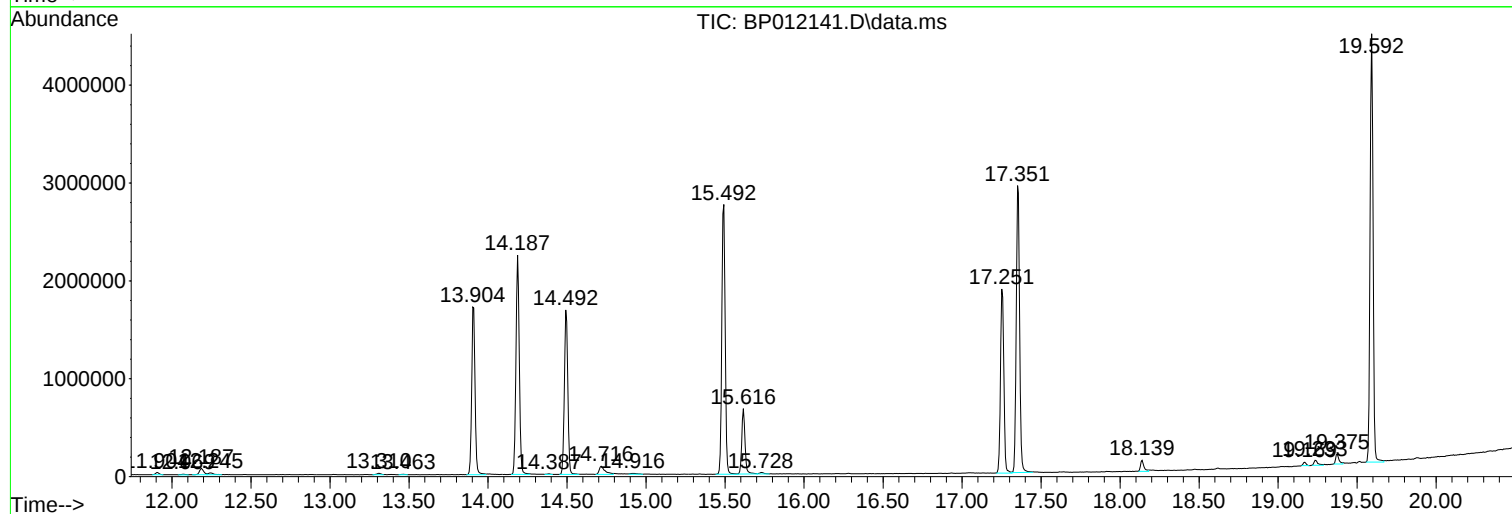
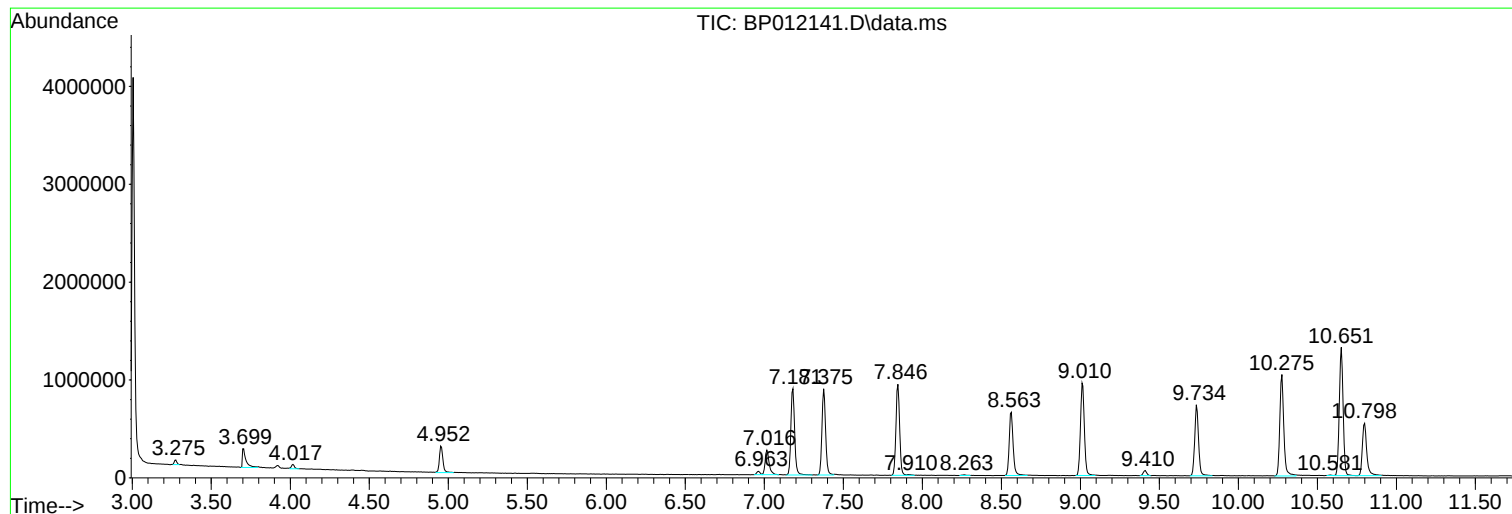
Sum of corrected areas: 56152513

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

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



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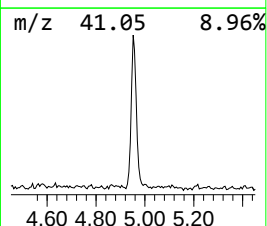
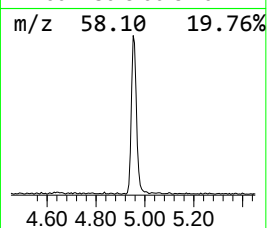
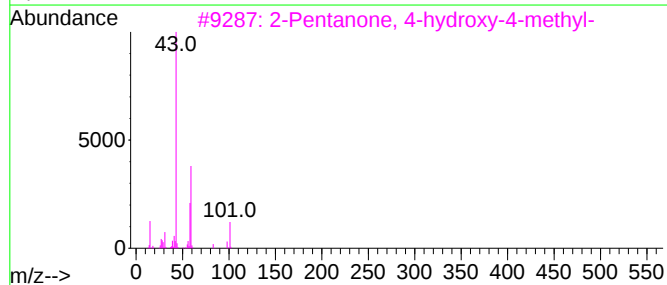
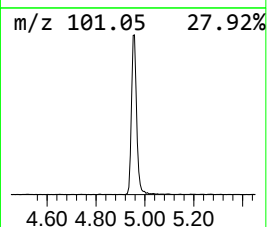
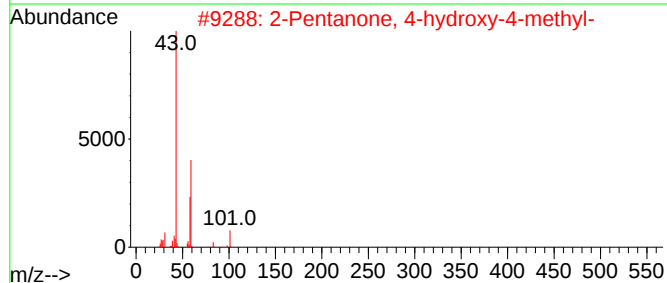
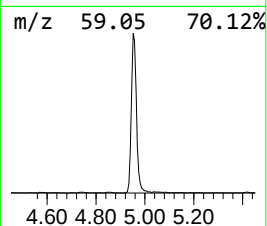
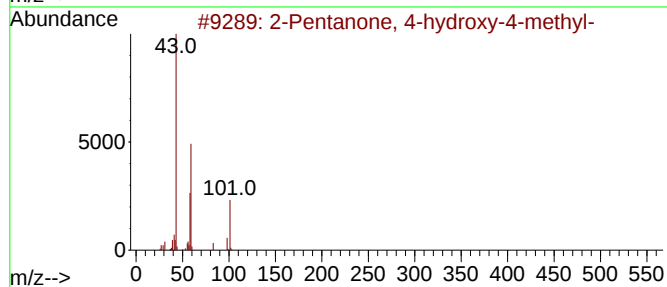
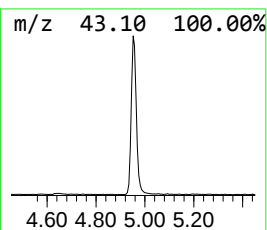
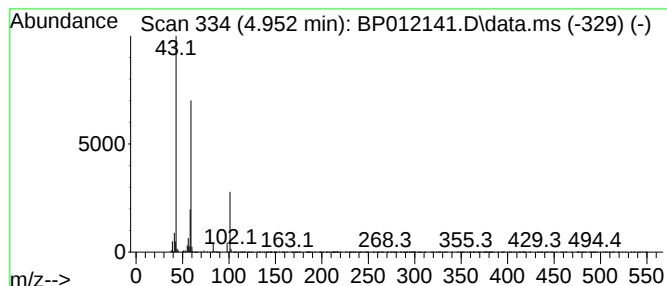
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Quant Title : SVOA CALIBRATION

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.
4.952	5.45 ng/ul	410264	1,4-Dichlorobenzene-d4	7.846

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	45		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42		
4	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	16		
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	12		



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Quant Title : SVOA CALIBRATION

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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-...	4.952	5.5	ng/u1	410264	1	7.846	1506040	20.0