

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051922\
 Data File : BP010410.D
 Acq On : 19 May 2022 19:30
 Operator : CG/JU
 Sample : N2918-06
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBTE4

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-BP051322.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010410.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.340	40	43	56	rVB	32821	60630	1.21%	0.142%
2	3.764	112	115	139	rVB	129906	268345	5.37%	0.629%
3	3.993	152	154	158	rVB2	9542	10216	0.20%	0.024%
4	4.087	166	170	174	rBV3	11163	15464	0.31%	0.036%
5	4.693	270	273	277	rBV3	7467	11544	0.23%	0.027%
6	5.011	321	327	342	rBV	1210374	1842465	36.87%	4.316%
7	6.463	569	574	576	rBV6	3247	5598	0.11%	0.013%
8	7.063	669	676	687	rVB	76942	139946	2.80%	0.328%
9	7.222	697	703	715	rBV	486327	824966	16.51%	1.932%
10	7.428	731	738	751	rBV	309529	522312	10.45%	1.223%
11	7.899	810	818	826	rBV	563899	955139	19.11%	2.237%
12	8.605	930	938	952	rBV	278020	487796	9.76%	1.143%
13	9.057	1007	1015	1026	rBV	636007	1146970	22.95%	2.687%
14	9.499	1085	1090	1095	rVB5	8761	13895	0.28%	0.033%
15	9.781	1130	1138	1150	rBV	331348	566664	11.34%	1.327%
16	10.110	1191	1194	1199	rBV6	3401	6045	0.12%	0.014%
17	10.322	1223	1230	1244	rBV	447105	820068	16.41%	1.921%
18	10.699	1286	1294	1300	rBV	822031	1399983	28.02%	3.279%
19	10.746	1300	1302	1309	rVB	23929	34279	0.69%	0.080%
20	10.834	1309	1317	1337	rBV	655180	1262823	25.27%	2.958%
21	11.522	1428	1434	1435	rBV5	3330	5216	0.10%	0.012%
22	12.134	1536	1538	1544	rVB6	4261	6633	0.13%	0.016%
23	12.234	1548	1555	1561	rBV2	26516	53962	1.08%	0.126%
24	12.293	1561	1565	1571	rVB4	11682	21578	0.43%	0.051%
25	12.587	1608	1615	1620	rVB4	2787	6807	0.14%	0.016%
26	13.269	1725	1731	1735	rBV9	3435	6688	0.13%	0.016%
27	13.363	1743	1747	1753	rVB3	13033	20211	0.40%	0.047%
28	13.510	1766	1772	1778	rVB3	2581	5253	0.11%	0.012%
29	13.940	1838	1845	1859	rBV	1640456	2375674	47.54%	5.565%
30	14.222	1882	1893	1908	rBV	1641579	2522198	50.47%	5.908%
31	14.434	1925	1929	1935	rVB4	8717	12828	0.26%	0.030%
32	14.528	1935	1945	1955	rBV	1281469	1961815	39.26%	4.595%
33	14.745	1975	1982	1996	rBV4	25411	84669	1.69%	0.198%
34	14.940	2011	2015	2023	rVB4	5157	11227	0.22%	0.026%
35	15.187	2050	2057	2063	rVV10	4469	8838	0.18%	0.021%
36	15.275	2067	2072	2078	rVV	21579	32195	0.64%	0.075%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
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37	15.345	2078	2084	2088	rVV7	8280	16793	0.34%	0.039%
38	15.381	2088	2090	2095	rVB3	6417	8222	0.16%	0.019%
39	15.522	2107	2114	2128	rBV	2337078	3462679	69.30%	8.111%
40	15.645	2129	2135	2147	rVV	441495	667329	13.35%	1.563%
41	15.763	2149	2155	2163	rVB3	29951	60164	1.20%	0.141%
42	15.892	2171	2177	2180	rVB8	3912	7664	0.15%	0.018%
43	16.251	2234	2238	2244	rVB7	8915	15030	0.30%	0.035%
44	16.310	2244	2248	2254	rVB7	13460	19889	0.40%	0.047%
45	16.428	2262	2268	2272	rBV9	4335	9012	0.18%	0.021%
46	16.786	2327	2329	2335	rVB7	6466	9175	0.18%	0.021%
47	17.045	2370	2373	2375	rBV3	8345	9751	0.20%	0.023%
48	17.210	2397	2401	2406	rBV8	5463	11643	0.23%	0.027%
49	17.281	2406	2413	2423	rBV	1599039	2371073	47.45%	5.554%
50	17.381	2423	2430	2446	rVV2	2866305	4249337	85.04%	9.953%
51	17.675	2474	2480	2482	rBV7	5175	9467	0.19%	0.022%
52	17.781	2496	2498	2507	rVB10	5234	8148	0.16%	0.019%
53	17.857	2509	2511	2516	rVB6	5777	6357	0.13%	0.015%
54	17.951	2523	2527	2532	rBV7	9338	13397	0.27%	0.031%
55	18.169	2559	2564	2573	rBV3	51472	91013	1.82%	0.213%
56	18.233	2573	2575	2582	rVB3	9561	17311	0.35%	0.041%
57	18.457	2610	2613	2617	rVB6	5284	6227	0.12%	0.015%
58	18.628	2640	2642	2645	rBV4	5363	5962	0.12%	0.014%
59	19.192	2734	2738	2745	rVB2	39396	53081	1.06%	0.124%
60	19.263	2745	2750	2754	rBV	36917	56997	1.14%	0.134%
61	19.398	2770	2773	2780	rBV6	23226	38555	0.77%	0.090%
62	19.545	2795	2798	2803	rVB6	17404	20635	0.41%	0.048%
63	19.616	2804	2810	2822	rBV	3725776	4996989	100.00%	11.705%
64	21.074	3054	3058	3064	rBV3	110622	179974	3.60%	0.422%
65	21.369	3102	3108	3123	rBV	1718503	2532564	50.68%	5.932%
66	23.639	3486	3494	3500	rBV2	2040748	4062550	81.30%	9.516%
67	23.792	3513	3520	3531	rVB2	1028791	2144729	42.92%	5.024%

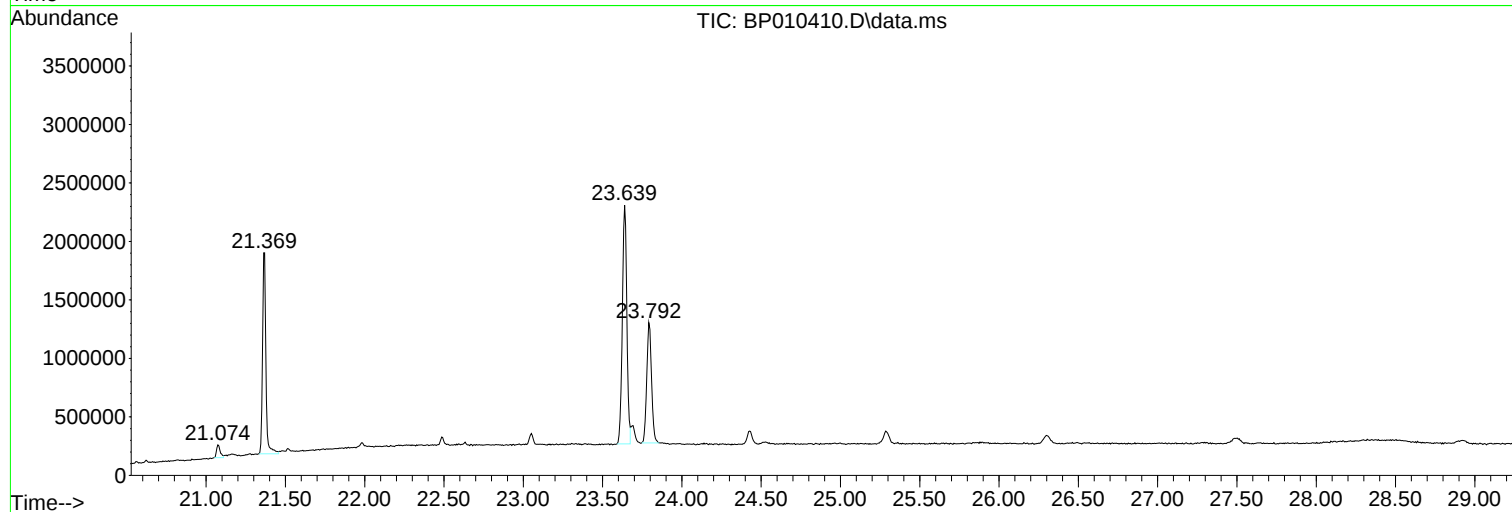
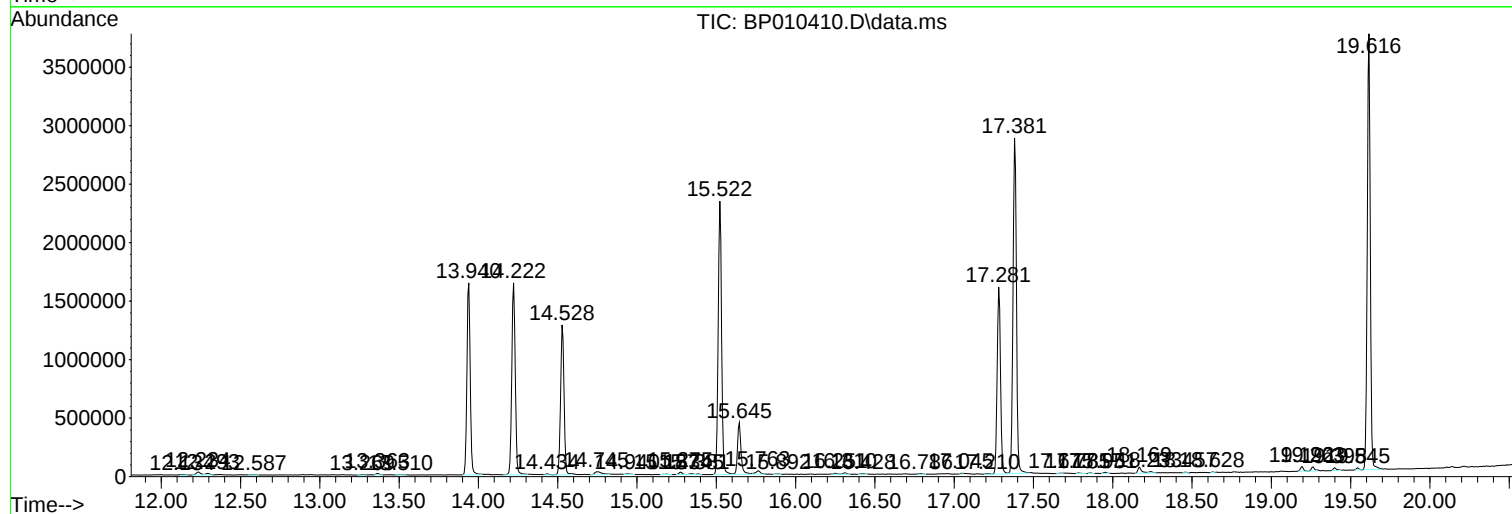
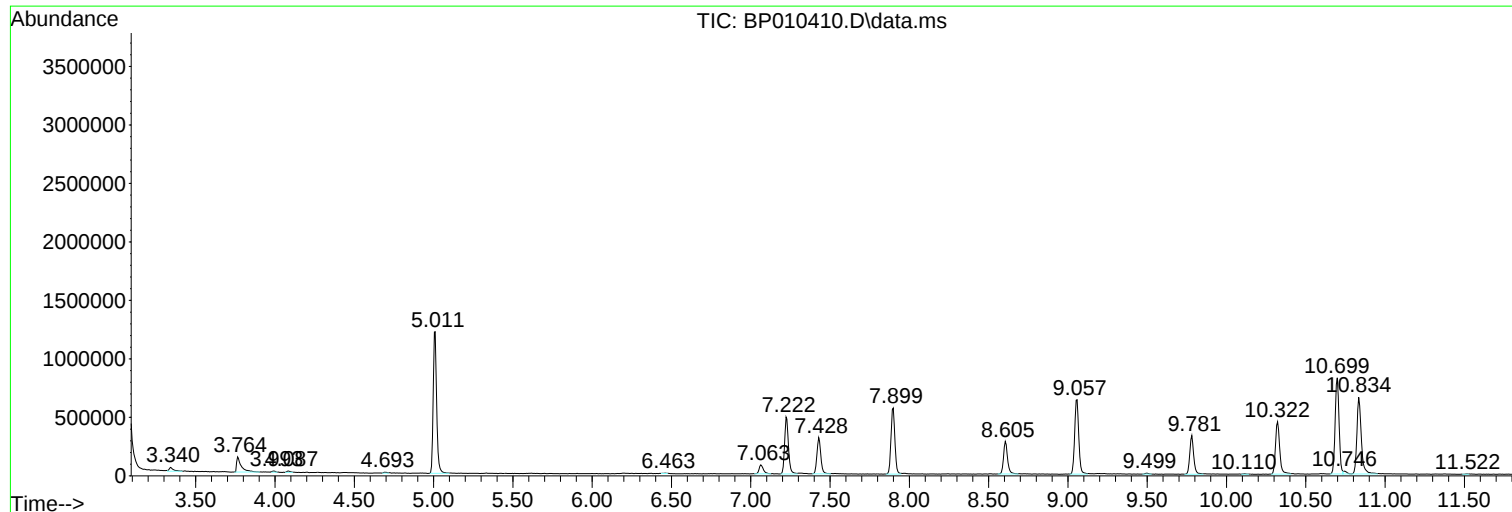
Sum of corrected areas: 42692657

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ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
BNA_P
ClientSampleId :
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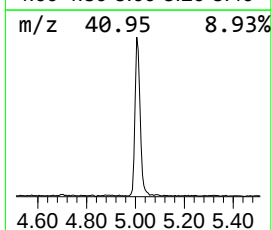
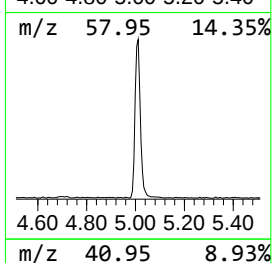
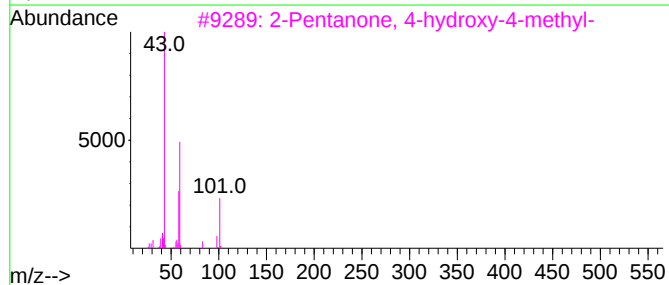
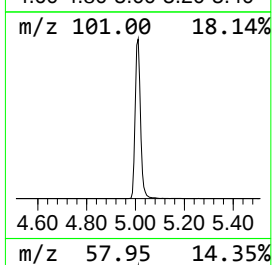
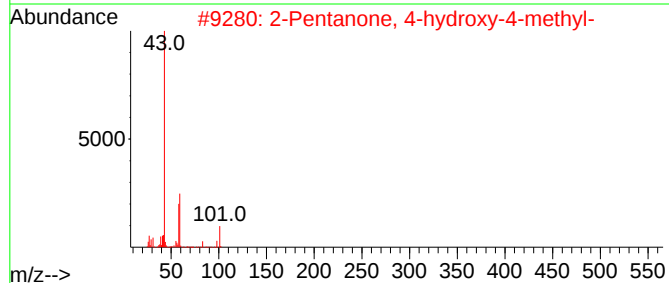
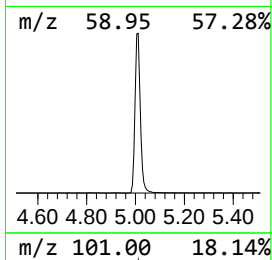
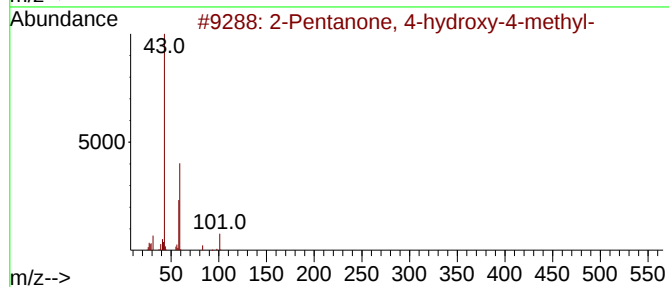
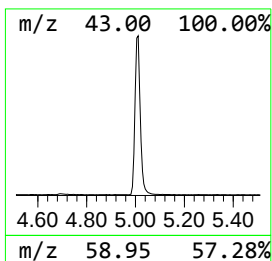
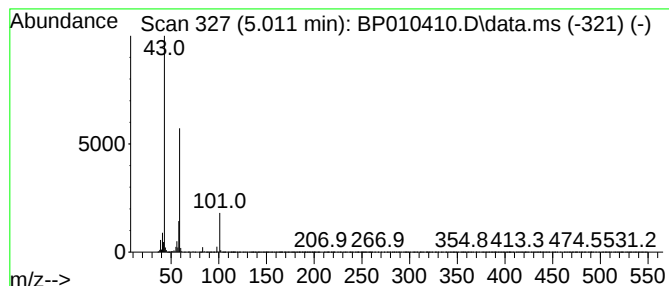
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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.
5.011	38.58 ng/ul	1842470	1,4-Dichlorobenzene-d4	7.899

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	40		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
5	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33		



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.011	38.6	ng/ul	1842470	1	7.899	955139	20.0