

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011979.D
 Acq On : 06 Oct 2022 20:19
 Operator : CG/JU
 Sample : N4923-18
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E10040

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-BP093022.M
 Title : SVOA CALIBRATION

Signal : TIC: BP011979.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.028	3	7	25	rVB	5301050	6560801	100.00%	11.293%
2	3.299	49	53	58	rVB	51226	56950	0.87%	0.098%
3	3.728	124	126	139	rBV	183922	315921	4.82%	0.544%
4	3.946	159	163	169	rVB	34458	49418	0.75%	0.085%
5	4.046	176	180	186	rVB	19073	25250	0.38%	0.043%
6	4.575	267	270	275	rVB6	6817	8313	0.13%	0.014%
7	4.981	334	339	349	rBV	146472	213479	3.25%	0.367%
8	5.157	365	369	372	rVB4	4305	7378	0.11%	0.013%
9	7.046	684	690	702	rBV	166361	290938	4.43%	0.501%
10	7.210	711	718	734	rBV	742955	1211237	18.46%	2.085%
11	7.410	744	752	763	rBV	672332	1114724	16.99%	1.919%
12	7.875	824	831	848	rVB	898150	1490652	22.72%	2.566%
13	8.592	945	953	964	rBV	413935	733802	11.18%	1.263%
14	9.045	1022	1030	1040	rBV	763102	1313840	20.03%	2.261%
15	9.769	1146	1153	1164	rBV	575781	969347	14.77%	1.668%
16	10.304	1236	1244	1258	rBV	780032	1436361	21.89%	2.472%
17	10.686	1301	1309	1323	rVB	1190482	2065040	31.48%	3.554%
18	10.828	1326	1333	1353	rBV	483800	977057	14.89%	1.682%
19	11.510	1442	1449	1458	rBV	29953	66615	1.02%	0.115%
20	12.222	1564	1570	1575	rBV	34138	70488	1.07%	0.121%
21	12.275	1576	1579	1589	rVB	18798	38190	0.58%	0.066%
22	13.422	1769	1774	1781	rBV9	7367	14057	0.21%	0.024%
23	13.933	1855	1861	1879	rBV	1687038	2416585	36.83%	4.159%
24	14.216	1901	1909	1924	rBV	2042686	2998344	45.70%	5.161%
25	14.522	1954	1961	1975	rBV	1614501	2473442	37.70%	4.257%
26	14.745	1991	1999	2018	rBV2	55180	198456	3.02%	0.342%
27	15.516	2124	2130	2136	rBV	2576166	3943059	60.10%	6.787%
28	15.563	2136	2138	2145	rVV	113371	171259	2.61%	0.295%
29	15.633	2145	2150	2167	rVV	629124	983552	14.99%	1.693%
30	15.757	2167	2171	2178	rVB3	12865	23295	0.36%	0.040%
31	17.280	2423	2430	2440	rBV	2000775	2846033	43.38%	4.899%
32	17.380	2440	2447	2462	rVV	2919108	4201182	64.03%	7.231%
33	18.163	2575	2580	2586	rBV4	38659	66678	1.02%	0.115%
34	19.192	2752	2755	2761	rVB3	32568	42281	0.64%	0.073%
35	19.262	2763	2767	2771	rBV2	43309	69452	1.06%	0.120%
36	19.615	2821	2827	2836	rBV	4214203	5404537	82.38%	9.302%

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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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Title : SVOA CALIBRATION

37	21.362	3120	3124	3131	rBV2	2619416	3526756	53.75%	6.070%
38	23.639	3504	3511	3526	rVB2	2949251	6023070	91.80%	10.367%
39	23.792	3531	3537	3547	rVB2	1812082	3680382	56.10%	6.335%

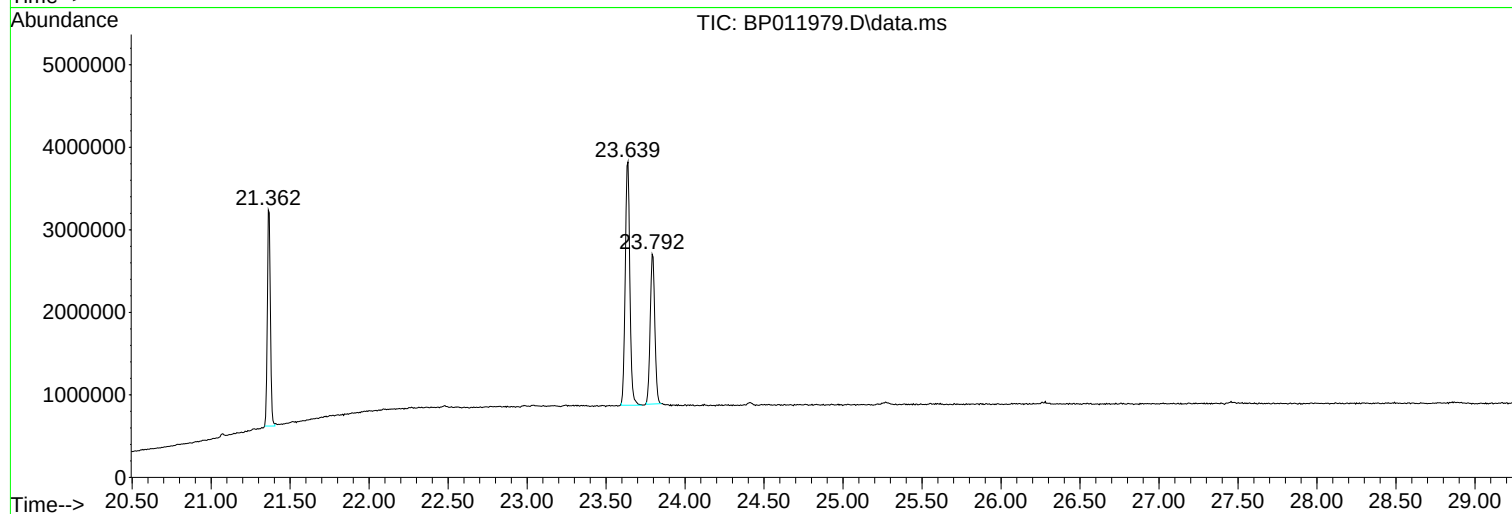
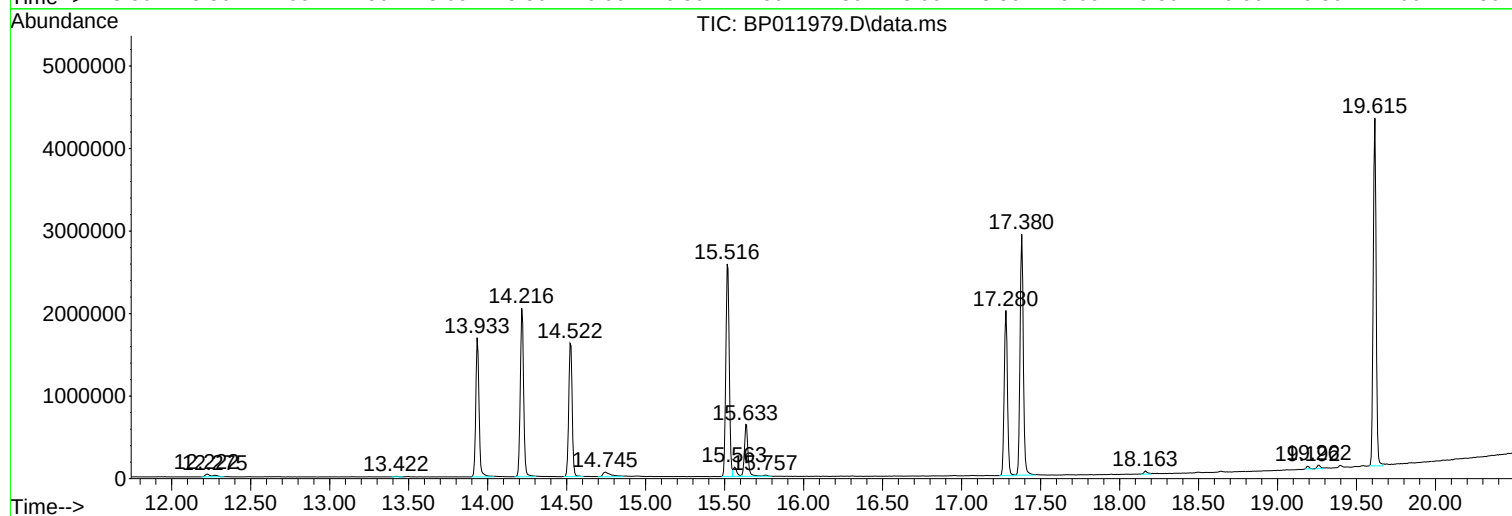
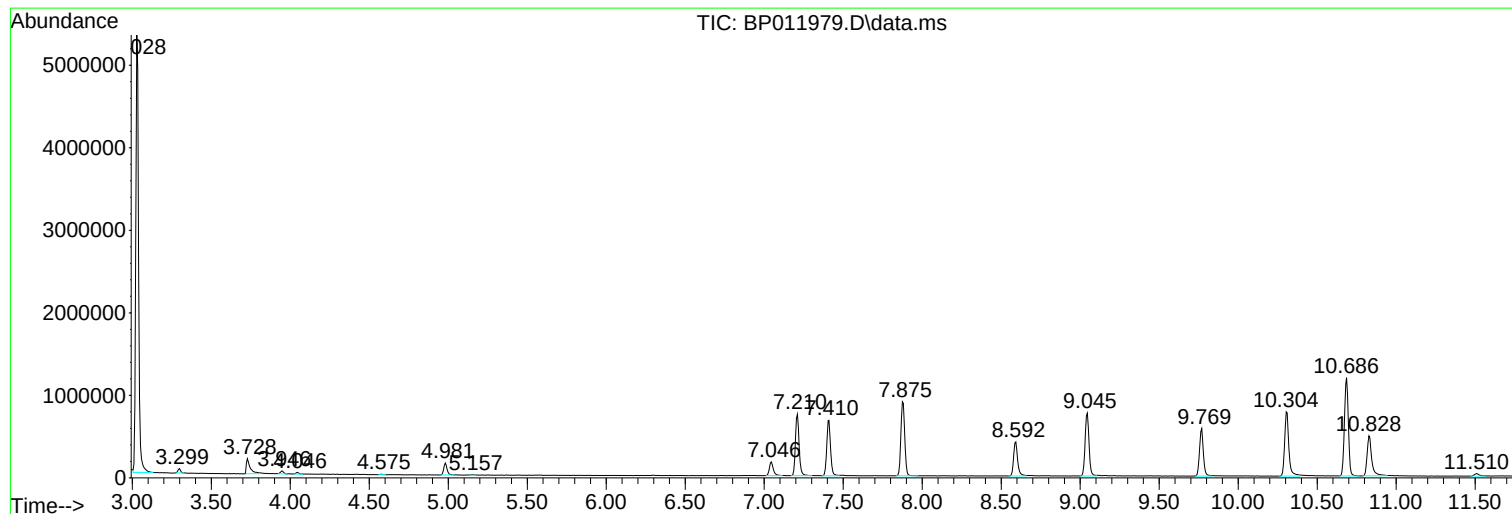
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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



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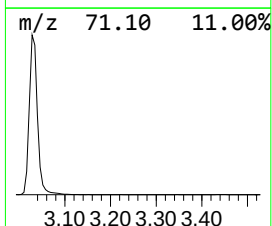
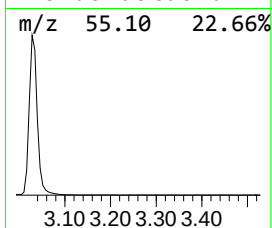
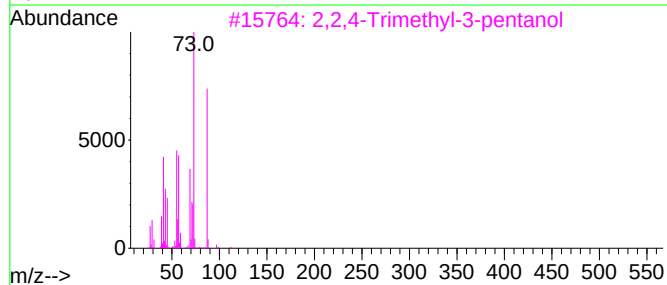
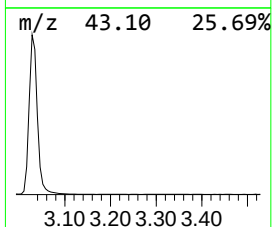
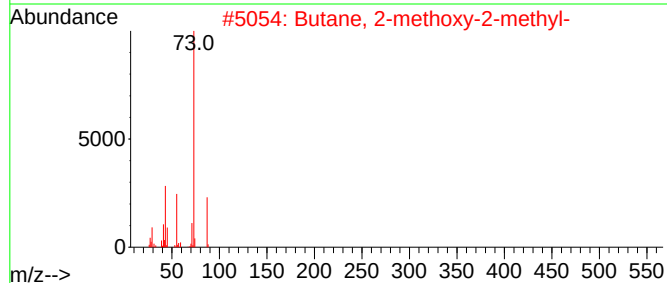
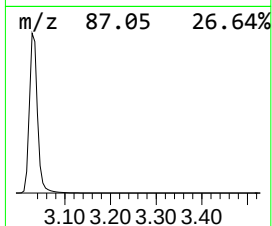
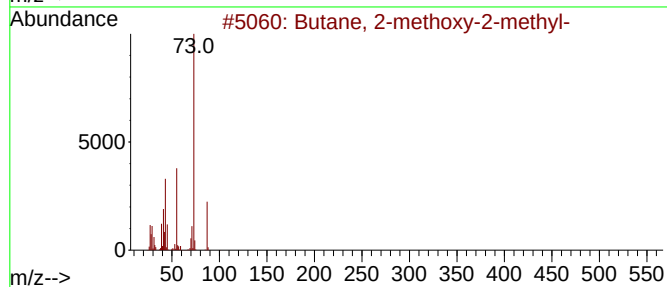
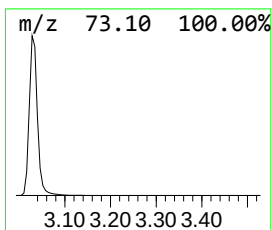
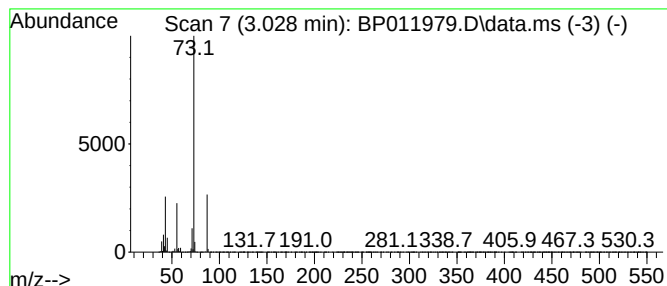
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.028	88.03 ng/ul	6560800	1,4-Dichlorobenzene-d4	7.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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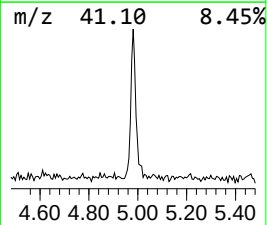
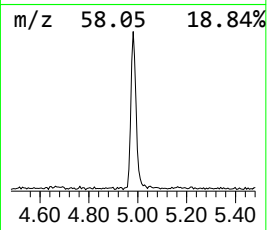
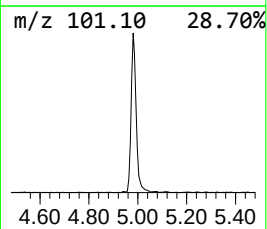
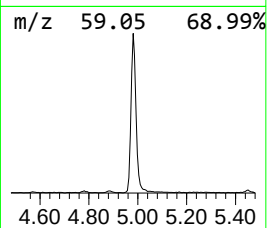
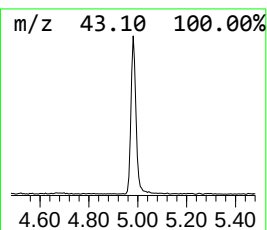
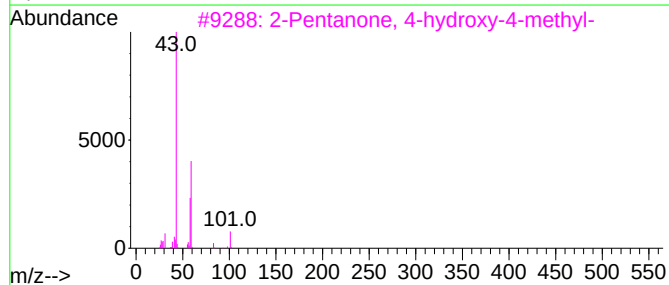
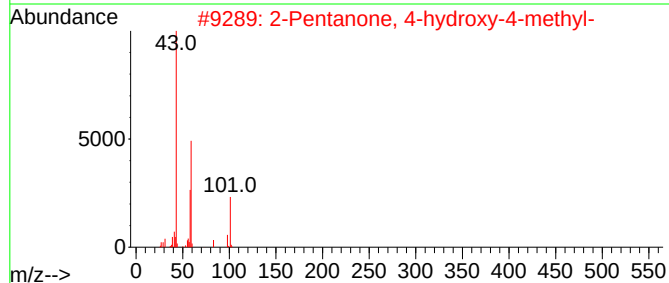
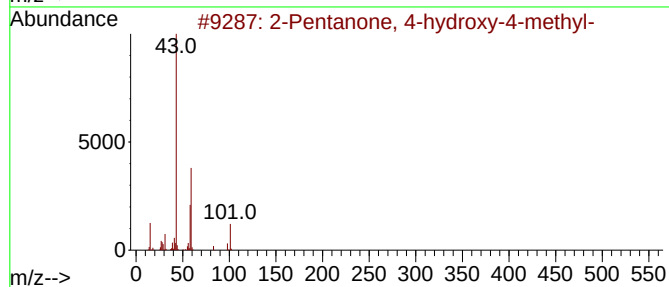
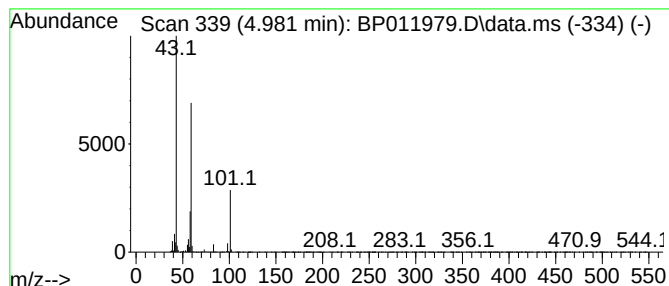
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.981	2.86 ng/ul	213479	1,4-Dichlorobenzene-d4	7.881

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	47		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45		
4	Guanidine	59	CH5N3	000113-00-8	43		
5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		



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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.028	88.0	ng/u1	6560800	1	7.881	1490650	20.0
2-Pentanone, 4-...	4.981	2.9	ng/u1	213479	1	7.881	1490650	20.0