

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP012007.D
 Acq On : 07 Oct 2022 13:47
 Operator : CG/JU
 Sample : PB148168BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK168

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: BP012007.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	15	rVB	120094	146800	3.17%	0.316%
2	3.293	49	52	57	rVB	47593	59676	1.29%	0.129%
3	3.723	123	125	148	rVB	513455	798869	17.25%	1.721%
4	7.040	683	689	704	rBV	699032	1163545	25.13%	2.506%
5	7.211	711	718	729	rVB	608218	995974	21.51%	2.145%
6	7.405	744	751	764	rBV	776718	1251488	27.03%	2.695%
7	7.875	823	831	843	rBV	680916	1093297	23.61%	2.355%
8	8.587	945	952	973	rBV	853535	1505277	32.51%	3.242%
9	9.046	1022	1030	1042	rVV	654660	1140722	24.64%	2.457%
10	9.763	1144	1152	1168	rBV	520617	932957	20.15%	2.009%
11	10.305	1237	1244	1263	rBV	803387	1488075	32.14%	3.205%
12	10.681	1301	1308	1320	rBV	919027	1596080	34.47%	3.438%
13	10.828	1326	1333	1350	rBV	739260	1433826	30.97%	3.088%
14	12.275	1574	1579	1588	rBV3	14006	27742	0.60%	0.060%
15	13.934	1853	1861	1878	rBV	1289653	1975862	42.67%	4.256%
16	14.216	1902	1909	1927	rBV2	1751359	2757865	59.56%	5.940%
17	14.522	1955	1961	1979	rBV	1358623	2077511	44.87%	4.474%
18	14.734	1990	1997	2020	rBV	335245	905355	19.55%	1.950%
19	15.204	2068	2077	2079	rBV10	2131	5524	0.12%	0.012%
20	15.522	2123	2131	2143	rBV	2165829	3439331	74.28%	7.408%
21	15.640	2146	2151	2167	rVV	415732	680212	14.69%	1.465%
22	15.757	2169	2171	2177	rVB4	14221	19864	0.43%	0.043%
23	16.092	2224	2228	2233	rBV8	3452	5823	0.13%	0.013%
24	17.069	2389	2394	2398	rBV5	4620	8245	0.18%	0.018%
25	17.281	2424	2430	2441	rBV	1629560	2435103	52.59%	5.245%
26	17.381	2441	2447	2462	rVB	2604203	3844151	83.03%	8.279%
27	19.192	2751	2755	2761	rBV2	32680	55014	1.19%	0.118%
28	19.263	2763	2767	2772	rBV	30542	49355	1.07%	0.106%
29	19.616	2821	2827	2842	rBV	3153722	4469522	96.53%	9.626%
30	21.369	3120	3125	3136	rBV	1972062	2752560	59.45%	5.928%
31	23.645	3505	3512	3529	rVB2	2214532	4630044	100.00%	9.972%
32	23.804	3533	3539	3551	rVB2	1277902	2684578	57.98%	5.782%

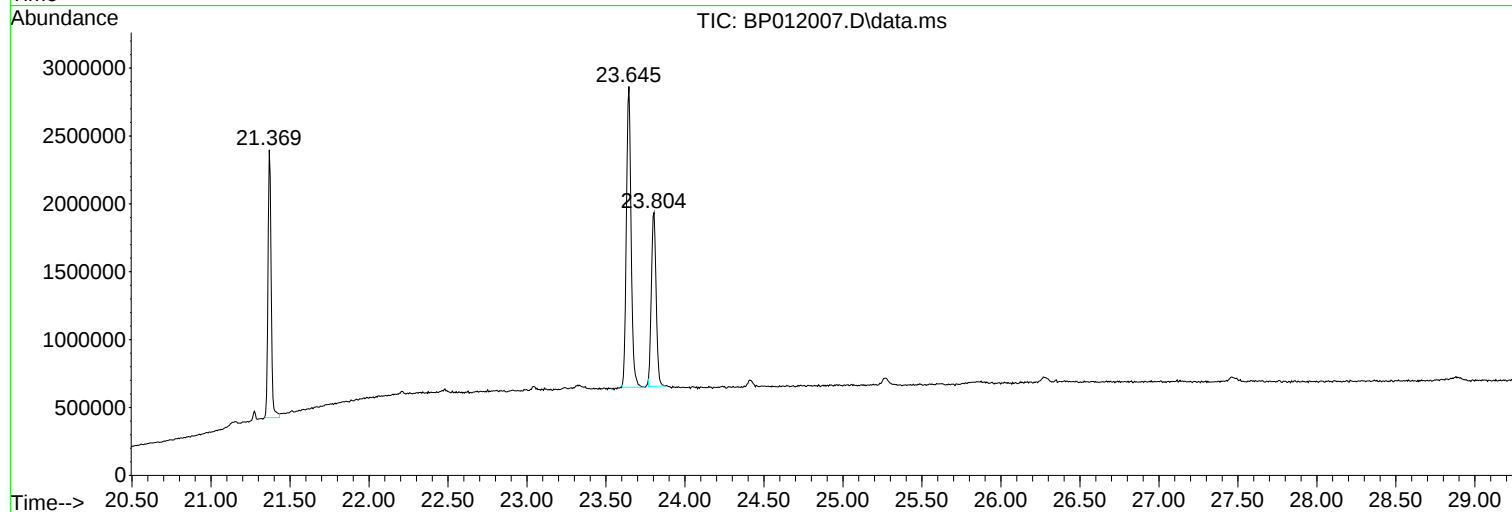
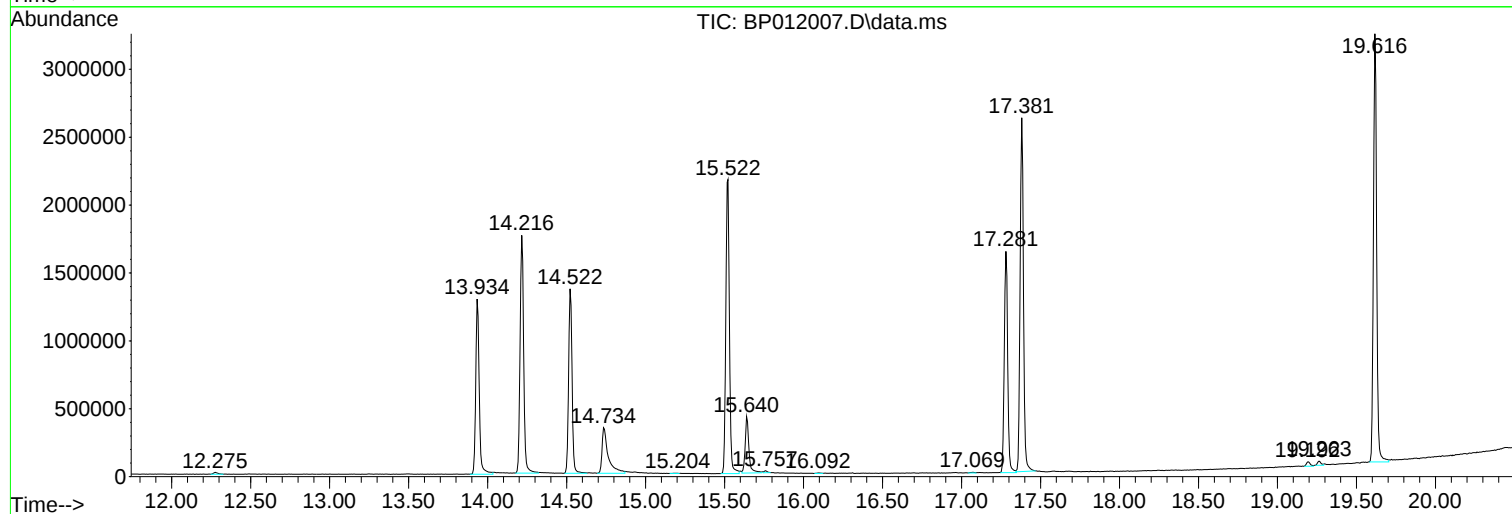
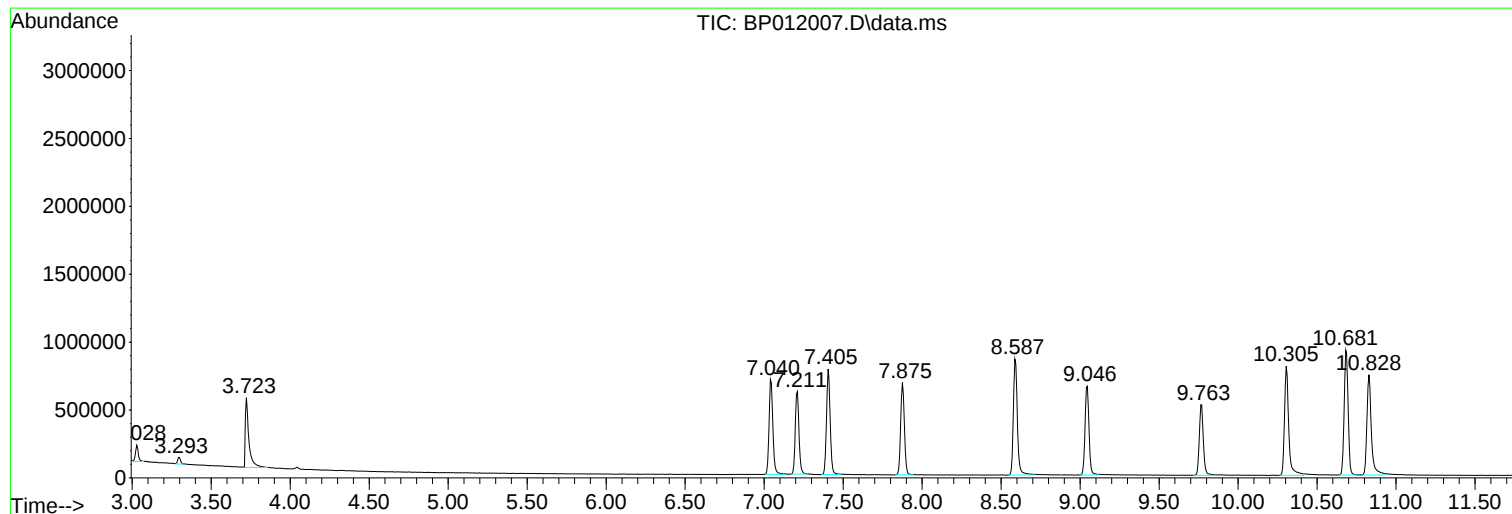
Sum of corrected areas: 46430247

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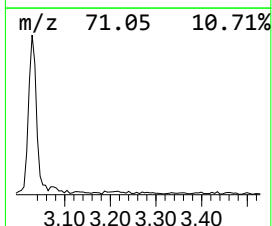
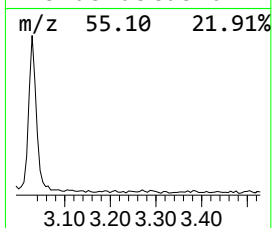
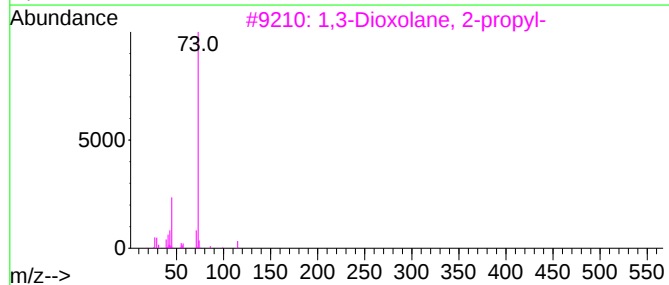
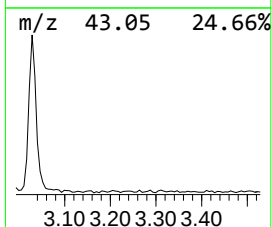
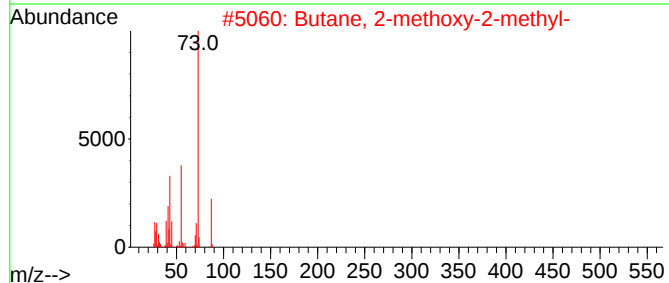
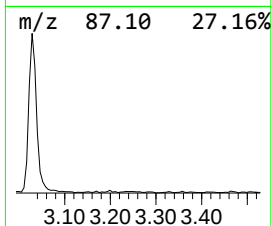
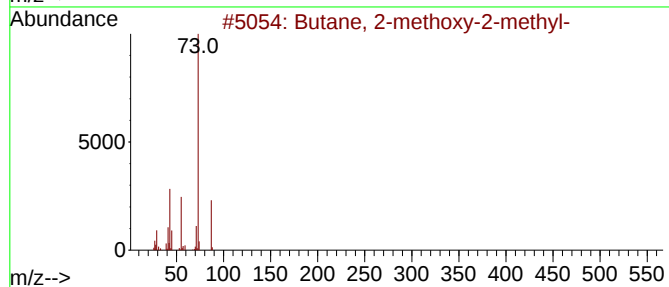
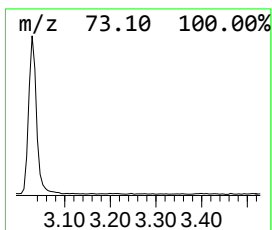
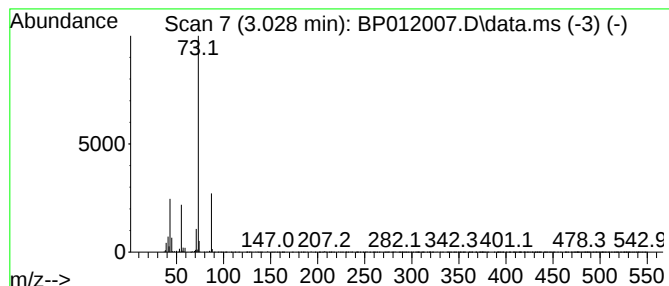
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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	2.69 ng/ul	146800	1,4-Dichlorobenzene-d4	7.875

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			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	25
4			N-Ethylformamide	73	C3H7NO	000627-45-2	9
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.028	2.7	ng/ul	146800	1	7.875	1093300	20.0