

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
 Data File : BP019165.D
 Acq On : 29 Dec 2023 14:41
 Operator : MA/JU
 Sample : 06021-08
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH7Z0

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-BP122723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019165.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.199	15	19	23	rBV	205571	277450	4.38%	0.588%
2	3.234	23	25	29	rVB	51326	54674	0.86%	0.116%
3	3.658	94	97	110	rBV	148939	228218	3.60%	0.484%
4	3.887	133	136	140	rVB3	8286	10280	0.16%	0.022%
5	3.975	148	151	157	rBV2	13094	18045	0.28%	0.038%
6	4.240	194	196	202	rVB3	8818	11190	0.18%	0.024%
7	4.422	222	227	236	rVB	121435	167224	2.64%	0.355%
8	4.799	287	291	297	rVB	56005	77597	1.22%	0.165%
9	4.899	305	308	312	rBV3	4802	6807	0.11%	0.014%
10	6.981	656	662	675	rVB	184844	307973	4.86%	0.653%
11	7.140	683	689	699	rBV	692379	1063234	16.78%	2.255%
12	7.334	715	722	732	rBV	664666	1029055	16.24%	2.182%
13	7.805	794	802	811	rBV	647442	1038673	16.39%	2.203%
14	8.522	917	924	938	rBV	500350	827387	13.06%	1.755%
15	8.963	993	999	1009	rBV	709058	1173066	18.51%	2.488%
16	9.369	1064	1068	1075	rVB4	9767	17768	0.28%	0.038%
17	9.687	1114	1122	1139	rBV	509698	835751	13.19%	1.772%
18	10.222	1206	1213	1234	rBV	871282	1489216	23.50%	3.158%
19	10.598	1267	1277	1290	rVB	829449	1386483	21.88%	2.940%
20	10.746	1293	1302	1315	rBV	644933	1146594	18.10%	2.431%
21	11.869	1490	1493	1497	rVB5	8250	11482	0.18%	0.024%
22	12.198	1543	1549	1557	rBV	14561	29904	0.47%	0.063%
23	13.345	1739	1744	1751	rBV3	10587	18514	0.29%	0.039%
24	13.863	1825	1832	1846	rBV	1693436	2293809	36.20%	4.864%
25	14.139	1872	1879	1890	rBV	1895938	2781917	43.91%	5.899%
26	14.445	1924	1931	1940	rBV	1233412	1790413	28.26%	3.797%
27	14.669	1962	1969	1985	rBV	80535	191779	3.03%	0.407%
28	15.198	2054	2059	2063	rBV	18418	26469	0.42%	0.056%
29	15.439	2093	2100	2108	rBV	2514101	3656901	57.72%	7.755%
30	15.569	2116	2122	2130	rBV	546362	750720	11.85%	1.592%
31	15.686	2138	2142	2145	rVB3	9789	13911	0.22%	0.029%
32	16.228	2229	2234	2239	rVB3	9672	14096	0.22%	0.030%
33	17.016	2359	2368	2376	rBV2	18921	39419	0.62%	0.084%
34	17.198	2393	2399	2408	rBV	1535180	2230983	35.21%	4.731%
35	17.304	2410	2417	2434	rVB	2818763	4146222	65.44%	8.792%
36	18.098	2545	2552	2560	rBV2	51476	84522	1.33%	0.179%

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Max Peaks: 100

Peak Location: TOP

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37	18.516	2619	2623	2625	rBV4	9548	11698	0.18%	0.025%
38	19.122	2722	2726	2730	rBV	43159	53049	0.84%	0.112%
39	19.192	2734	2738	2742	rBV	37818	54167	0.85%	0.115%
40	19.339	2759	2763	2765	rBV4	28700	36680	0.58%	0.078%
41	19.480	2785	2787	2791	rVB4	13083	13137	0.21%	0.028%
42	19.545	2792	2798	2812	rBV	3966863	5320727	83.98%	11.283%
43	21.304	3092	3097	3105	rBV	2281280	2884442	45.52%	6.117%
44	23.515	3466	3473	3488	rVB2	3172254	6336051	100.00%	13.436%
45	23.668	3493	3499	3511	rVB	1596117	3199278	50.49%	6.784%

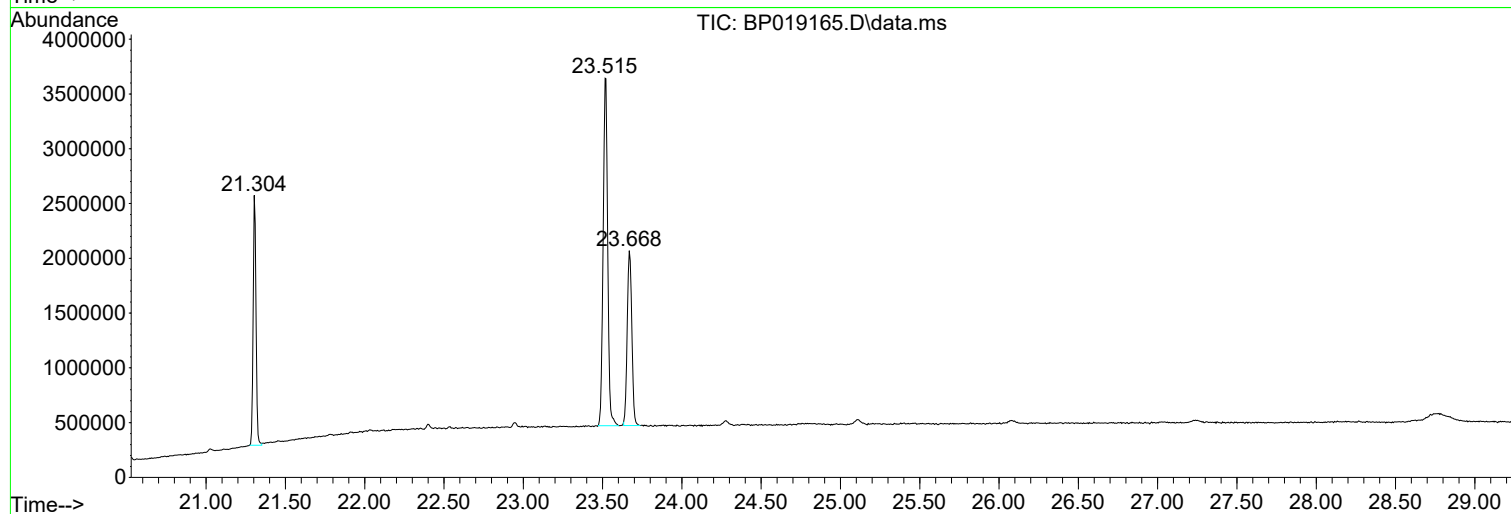
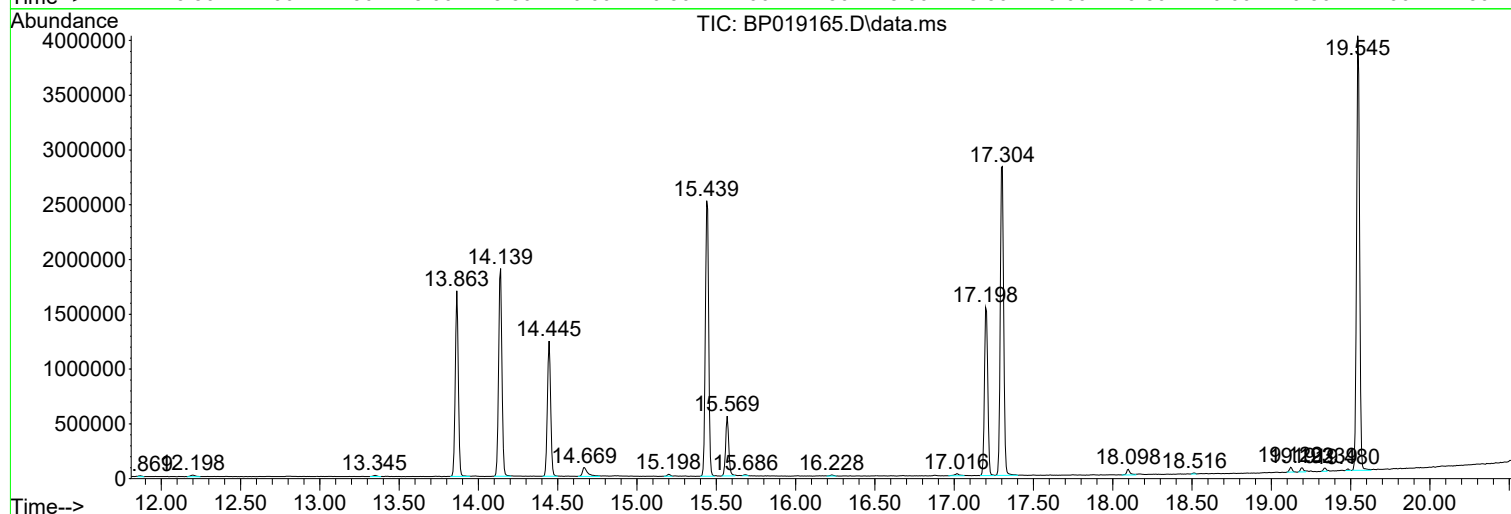
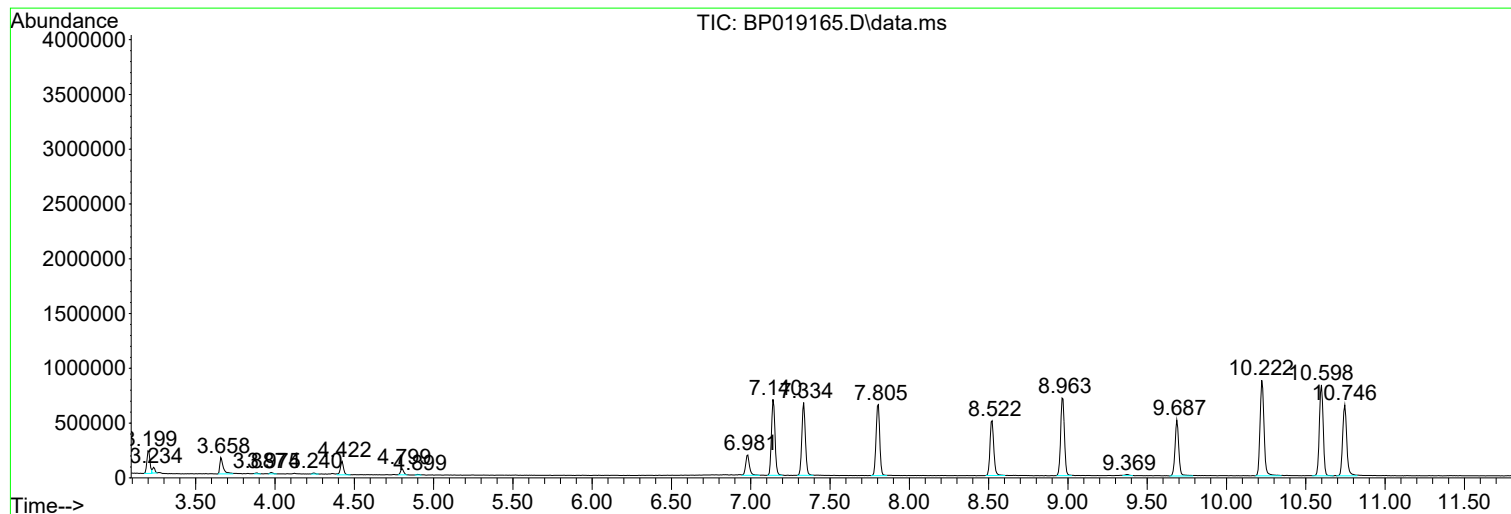
Sum of corrected areas: 47156975

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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



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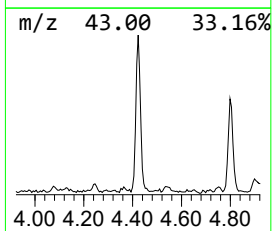
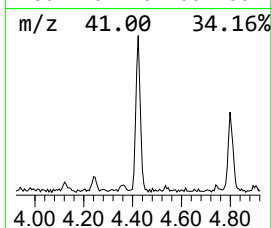
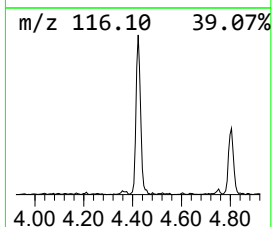
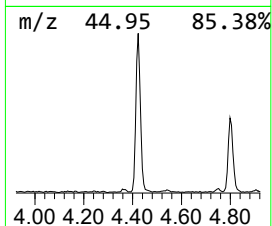
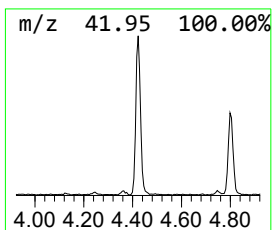
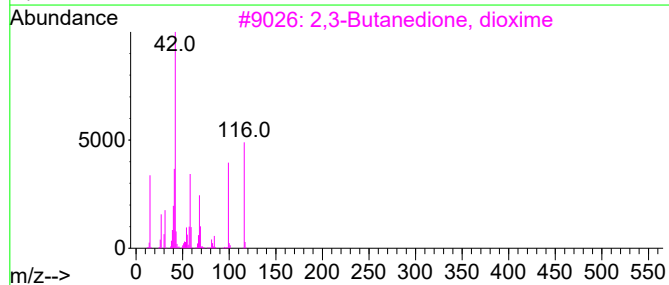
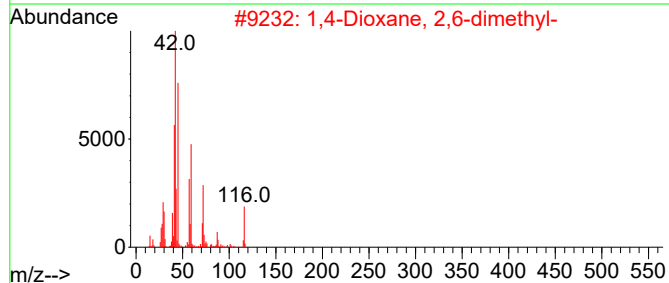
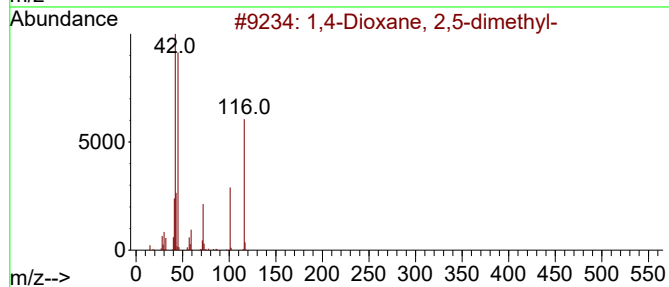
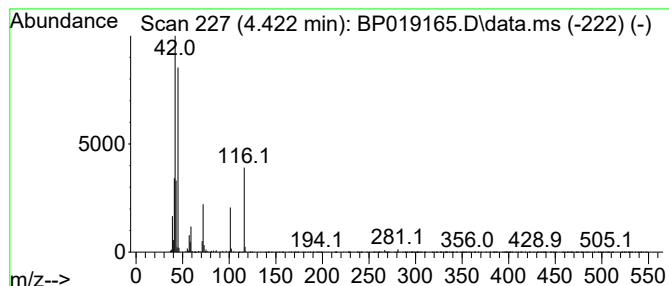
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 1,4-Dioxane, 2,5-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.422	3.22 ng/ul	167224	1,4-Dichlorobenzene-d4	7.805

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Dioxane, 2,5-dimethyl-	116	C6H12O2	015176-21-3	94		
2	1,4-Dioxane, 2,6-dimethyl-	116	C6H12O2	010138-17-7	38		
3	2,3-Butanedione, dioxime	116	C4H8N2O2	000095-45-4	23		
4	Ethyl hexan-2-yl carbonate	174	C9H18O3	1010373-79-4	16		
5	2-Butanol	74	C4H10O	000078-92-2	12		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1,4-Dioxane, 2,...	4.422	3.2	ng/ul	167224	1	7.805	1038670	20.0