

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009770.D
 Acq On : 12 Apr 2022 02:04
 Operator : CG/JU
 Sample : N2338-16
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J88

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-BP040722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP009770.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.128	3	7	24	rVB	291895	467153	9.77%	1.280%
2	3.387	47	51	59	rBV	23841	33941	0.71%	0.093%
3	3.675	94	100	112	rVB3	15109	39218	0.82%	0.107%
4	3.811	120	123	135	rBV	58278	109507	2.29%	0.300%
5	4.140	175	179	183	rVB5	4616	7402	0.15%	0.020%
6	5.075	334	338	342	rBV5	3370	4909	0.10%	0.013%
7	7.122	679	686	696	rVB	105801	177152	3.71%	0.485%
8	7.293	705	715	726	rBV	385835	615941	12.88%	1.688%
9	7.499	741	750	762	rBV	363764	599670	12.54%	1.643%
10	7.969	822	830	839	rBV	364571	594686	12.44%	1.630%
11	8.669	941	949	960	rBV	313708	516633	10.81%	1.416%
12	9.122	1018	1026	1039	rBV	517467	882944	18.47%	2.420%
13	9.240	1042	1046	1052	rVB8	2488	4932	0.10%	0.014%
14	9.463	1078	1084	1091	rBV7	10856	19970	0.42%	0.055%
15	9.587	1099	1105	1110	rBV6	7187	12793	0.27%	0.035%
16	9.851	1142	1150	1168	rVB	388003	644103	13.47%	1.765%
17	10.393	1234	1242	1256	rBV	543174	918898	19.22%	2.518%
18	10.569	1267	1272	1277	rBV9	3185	6572	0.14%	0.018%
19	10.775	1297	1307	1318	rBV	546515	931109	19.48%	2.552%
20	10.904	1322	1329	1339	rVV	415887	716013	14.98%	1.962%
21	12.087	1524	1530	1537	rBV5	4329	8024	0.17%	0.022%
22	12.357	1569	1576	1583	rBV2	14003	23278	0.49%	0.064%
23	13.445	1754	1761	1768	rBV8	3774	8382	0.18%	0.023%
24	14.004	1848	1856	1863	rBV	1426352	1954818	40.89%	5.357%
25	14.292	1894	1905	1920	rBV	1457306	2099572	43.92%	5.754%
26	14.422	1922	1927	1929	rBV5	3719	5959	0.12%	0.016%
27	14.598	1950	1957	1968	rBV	923178	1377618	28.82%	3.775%
28	14.775	1980	1987	1993	rBV	150550	234860	4.91%	0.644%
29	15.239	2061	2066	2071	rBV9	3933	6690	0.14%	0.018%
30	15.410	2092	2095	2099	rVB4	5245	7046	0.15%	0.019%
31	15.592	2119	2126	2136	rBV	1973976	2863059	59.89%	7.846%
32	15.698	2137	2144	2158	rVV	644205	902411	18.88%	2.473%
33	15.834	2161	2167	2173	rVB2	23703	36087	0.75%	0.099%
34	16.045	2194	2203	2211	rBV3	16493	29292	0.61%	0.080%
35	16.157	2217	2222	2227	rBV8	3961	6870	0.14%	0.019%
36	16.281	2239	2243	2247	rBV6	3233	5732	0.12%	0.016%

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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

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
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37	16.381	2256	2260	2266	rVB6	8327	13426	0.28%	0.037%
38	16.804	2326	2332	2336	rBV3	21784	32087	0.67%	0.088%
39	16.951	2352	2357	2364	rBV4	12049	20389	0.43%	0.056%
40	17.022	2366	2369	2375	rVB8	4683	7764	0.16%	0.021%
41	17.128	2383	2387	2391	rBV4	7062	11499	0.24%	0.032%
42	17.281	2410	2413	2416	rVB5	5726	6145	0.13%	0.017%
43	17.351	2417	2425	2434	rBV	1253559	1768945	37.00%	4.848%
44	17.451	2435	2442	2455	rBV	2477740	3581309	74.92%	9.815%
45	17.586	2461	2465	2470	rVB8	4760	7284	0.15%	0.020%
46	17.739	2489	2491	2496	rVB5	5214	6146	0.13%	0.017%
47	17.869	2508	2513	2516	rBV7	3443	6559	0.14%	0.018%
48	18.028	2535	2540	2546	rVB	91860	116702	2.44%	0.320%
49	18.233	2570	2575	2581	rBV4	22921	35123	0.73%	0.096%
50	18.533	2623	2626	2633	rVB9	8932	16349	0.34%	0.045%
51	18.651	2644	2646	2650	rBV4	4641	5064	0.11%	0.014%
52	19.151	2727	2731	2736	rVB	101713	111920	2.34%	0.307%
53	19.251	2744	2748	2752	rBV3	35772	53939	1.13%	0.148%
54	19.322	2756	2760	2764	rVV	35483	47613	1.00%	0.130%
55	19.675	2814	2820	2833	rBV	3345115	4398756	92.01%	12.055%
56	21.139	3066	3069	3073	rVB4	61740	81232	1.70%	0.223%
57	21.427	3113	3118	3125	rVB	1741501	2179123	45.58%	5.972%
58	23.739	3503	3511	3520	rVB	2375685	4780483	100.00%	13.101%
59	23.898	3531	3538	3551	rVB2	1107623	2328588	48.71%	6.381%

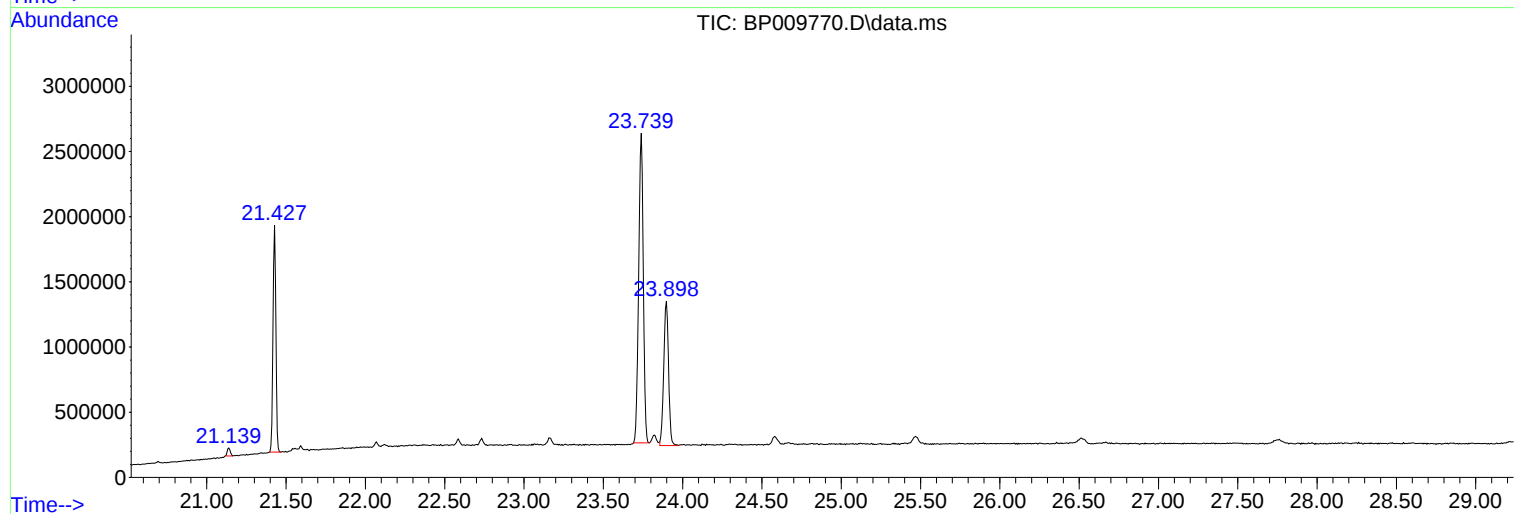
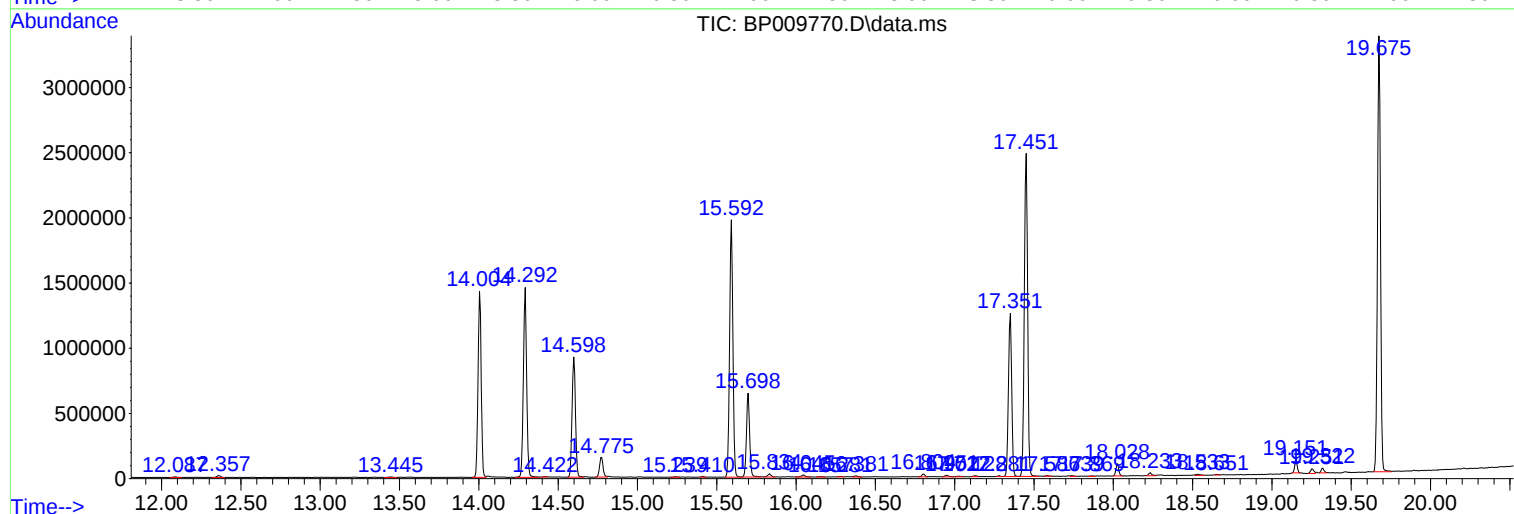
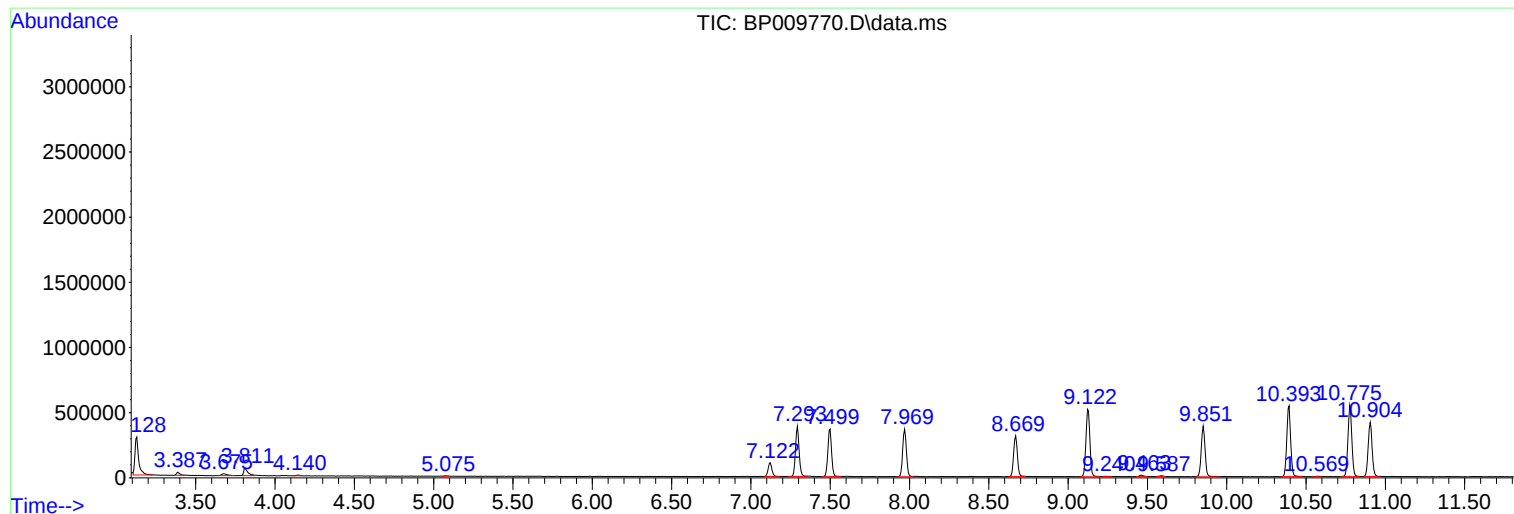
Sum of corrected areas: 36489689

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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



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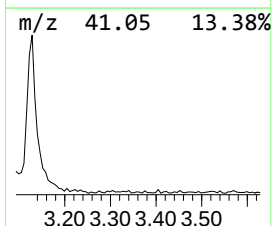
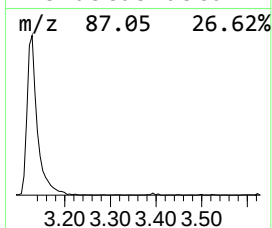
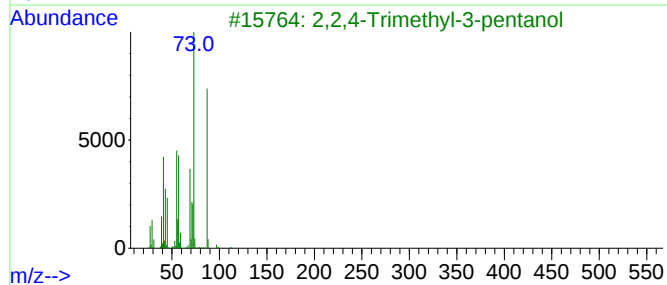
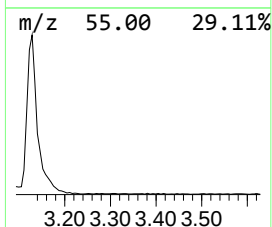
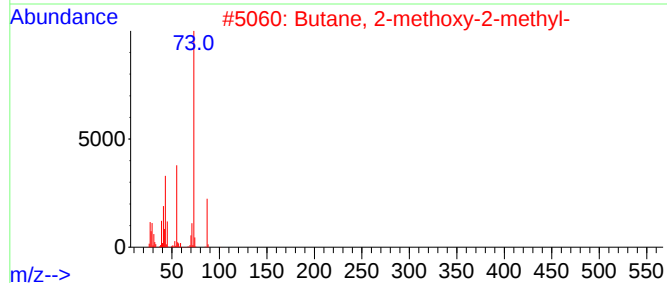
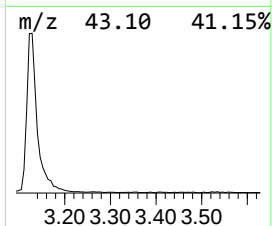
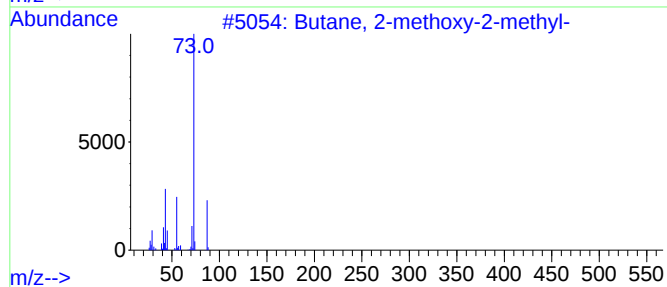
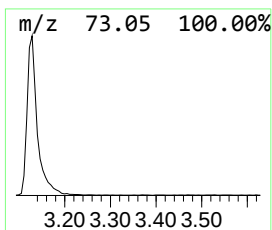
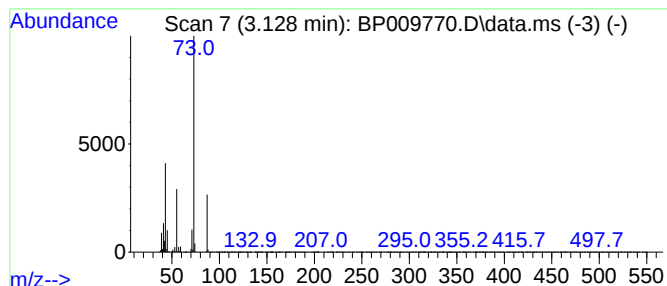
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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.
3.128	15.71 ng/ul	467153	1,4-Dichlorobenzene-d4	7.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	32
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	23



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.128	15.7	ng/ul	467153	1	7.969	594686	20.0