

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101422\
 Data File : BP012158.D
 Acq On : 15 Oct 2022 03:18
 Operator : CG/JU
 Sample : N4984-15
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGNC1

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-BP101122.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012158.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.011	678	684	697	rVB	628127	1061665	23.69%	2.348%
2	7.175	706	712	727	rVB	534266	866088	19.33%	1.916%
3	7.369	739	745	756	rBV	639992	1021894	22.80%	2.260%
4	7.840	819	825	833	rBV	790123	1256210	28.03%	2.778%
5	8.558	940	947	961	rBV	698112	1222429	27.28%	2.704%
6	8.675	963	967	976	rVB3	18037	34934	0.78%	0.077%
7	9.010	1017	1024	1038	rBV	568857	953204	21.27%	2.108%
8	9.163	1048	1050	1057	rVB7	8196	11752	0.26%	0.026%
9	9.405	1086	1091	1096	rVB2	9057	15681	0.35%	0.035%
10	9.728	1137	1146	1157	rBV	415170	721795	16.11%	1.596%
11	10.269	1229	1238	1258	rBV	681329	1194913	26.67%	2.643%
12	10.428	1258	1265	1281	rBV	338449	665439	14.85%	1.472%
13	10.646	1294	1302	1309	rBV	1069297	1788752	39.92%	3.956%
14	10.793	1320	1327	1350	rVB	550351	1036953	23.14%	2.293%
15	11.481	1439	1444	1449	rBV9	3635	8739	0.20%	0.019%
16	11.663	1470	1475	1481	rBV10	2389	6458	0.14%	0.014%
17	12.069	1539	1544	1547	rVB7	4584	6793	0.15%	0.015%
18	12.240	1568	1573	1581	rBV4	10086	19565	0.44%	0.043%
19	12.663	1641	1645	1651	rBV8	2500	5593	0.12%	0.012%
20	12.781	1661	1665	1668	rBV6	3165	6276	0.14%	0.014%
21	12.846	1670	1676	1678	rVV7	3156	5101	0.11%	0.011%
22	13.093	1713	1718	1725	rVB8	5890	11368	0.25%	0.025%
23	13.304	1751	1754	1761	rVB3	11278	15533	0.35%	0.034%
24	13.381	1761	1767	1773	rVV2	23017	37651	0.84%	0.083%
25	13.451	1773	1779	1784	rVB9	4364	8182	0.18%	0.018%
26	13.610	1802	1806	1811	rBV8	2472	4782	0.11%	0.011%
27	13.704	1816	1822	1828	rVV6	9092	16664	0.37%	0.037%
28	13.816	1834	1841	1845	rVB5	9087	13204	0.29%	0.029%
29	13.904	1849	1856	1870	rBV	1136416	1610540	35.94%	3.562%
30	14.140	1892	1896	1897	rBV4	5206	6169	0.14%	0.014%
31	14.181	1897	1903	1919	rVV	1423600	2148979	47.96%	4.753%
32	14.304	1922	1924	1928	rVV4	3349	5537	0.12%	0.012%
33	14.381	1928	1937	1943	rVB7	6572	12972	0.29%	0.029%
34	14.493	1948	1956	1963	rBV	1466121	2207215	49.26%	4.882%
35	14.551	1963	1966	1974	rVB	17103	29566	0.66%	0.065%
36	14.657	1980	1984	1986	rBV5	5695	7985	0.18%	0.018%

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Stop Thrs : 0

Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
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37	14.704	1986	1992	2013	rVV	300448	617729	13.78%	1.366%
38	14.916	2019	2028	2034	rBV2	50151	91434	2.04%	0.202%
39	15.004	2041	2043	2048	rVB6	3832	5525	0.12%	0.012%
40	15.157	2067	2069	2073	rVB5	3859	5088	0.11%	0.011%
41	15.287	2082	2091	2094	rBV8	6313	14400	0.32%	0.032%
42	15.340	2095	2100	2102	rVB6	4993	8777	0.20%	0.019%
43	15.487	2117	2125	2132	rBV	1953634	2708358	60.44%	5.990%
44	15.610	2141	2146	2154	rBV	209741	321376	7.17%	0.711%
45	15.722	2163	2165	2173	rVB8	9255	17417	0.39%	0.039%
46	15.840	2180	2185	2192	rVB5	10126	17166	0.38%	0.038%
47	15.945	2199	2203	2207	rBV5	9505	17169	0.38%	0.038%
48	15.987	2207	2210	2216	rVB	8759	13869	0.31%	0.031%
49	16.216	2243	2249	2254	rVB10	8035	17583	0.39%	0.039%
50	16.269	2254	2258	2264	rBV8	4950	7829	0.17%	0.017%
51	16.387	2275	2278	2283	rVB7	4623	7330	0.16%	0.016%
52	16.551	2297	2306	2312	rBV8	9754	29511	0.66%	0.065%
53	16.604	2312	2315	2318	rVB5	4356	5518	0.12%	0.012%
54	16.787	2342	2346	2349	rBV5	8284	11944	0.27%	0.026%
55	16.975	2374	2378	2379	rBV4	6252	8482	0.19%	0.019%
56	17.063	2390	2393	2400	rVB4	10428	18402	0.41%	0.041%
57	17.251	2418	2425	2429	rBV	1852986	2632861	58.75%	5.823%
58	17.292	2429	2432	2436	rVV	123989	173463	3.87%	0.384%
59	17.351	2436	2442	2454	rVV	2040180	3092933	69.02%	6.841%
60	17.451	2454	2459	2466	rVB6	31902	57009	1.27%	0.126%
61	17.534	2471	2473	2476	rBV3	4892	6407	0.14%	0.014%
62	17.916	2534	2538	2544	rBV	41332	50343	1.12%	0.111%
63	18.134	2568	2575	2578	rBV2	128262	202643	4.52%	0.448%
64	18.245	2591	2594	2599	rVB	22177	30738	0.69%	0.068%
65	18.310	2599	2605	2608	rBV4	56889	95460	2.13%	0.211%
66	18.604	2652	2655	2661	rVB2	25298	35130	0.78%	0.078%
67	19.004	2720	2723	2727	rBV4	15800	29989	0.67%	0.066%
68	19.051	2728	2731	2736	rVB6	28696	43290	0.97%	0.096%
69	19.163	2747	2750	2756	rVB4	30865	48358	1.08%	0.107%
70	19.263	2762	2767	2773	rVB	470178	622450	13.89%	1.377%
71	19.369	2781	2785	2788	rBV2	86913	115839	2.59%	0.256%
72	19.586	2816	2822	2835	rBV2	3013922	4481195	100.00%	9.911%
73	20.098	2906	2909	2913	rVB2	98350	123697	2.76%	0.274%
74	20.192	2922	2925	2930	rBV	80478	109149	2.44%	0.241%
75	21.039	3067	3069	3072	rBV2	106207	129903	2.90%	0.287%
76	21.339	3114	3120	3124	rBV	2676923	3683423	82.20%	8.147%

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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 >
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77	23.598	3498	3504	3510	rBV2	1939366	3847258	85.85%	8.509%
78	23.757	3524	3531	3546	rVB2	1731752	3611692	80.60%	7.988%

Sum of corrected areas: 45214718

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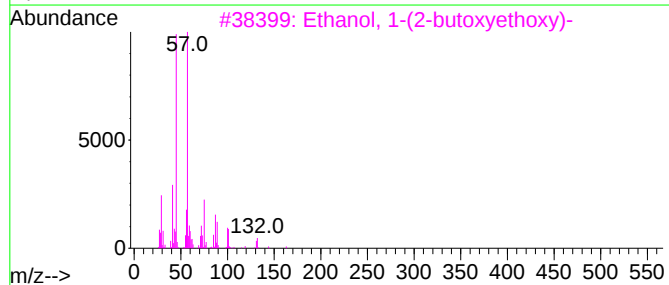
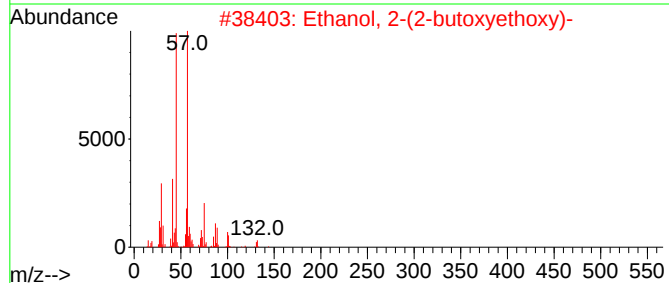
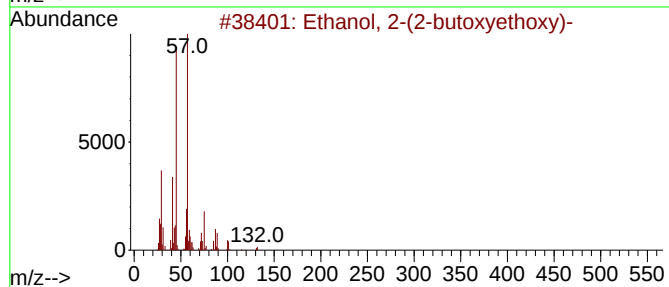
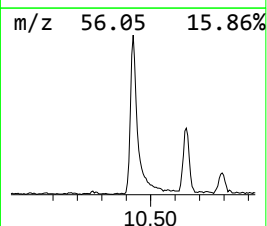
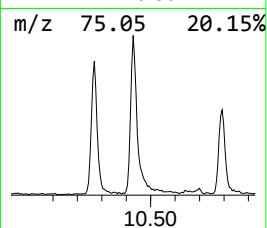
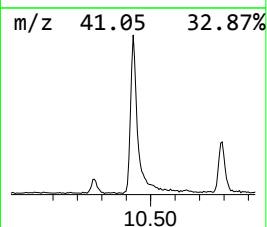
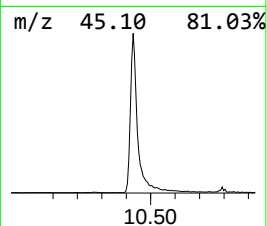
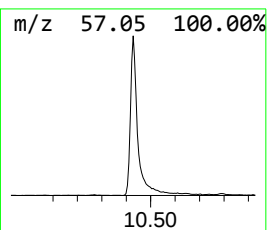
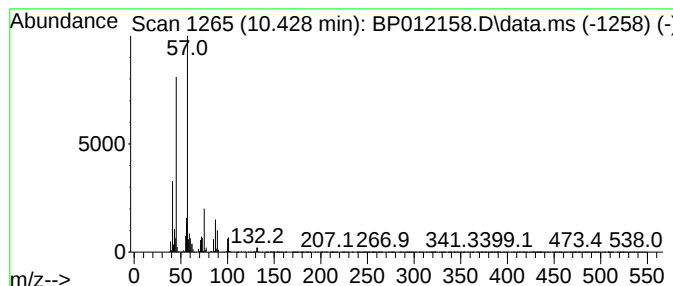
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.428	7.44 ng/ul	665439	Naphthalene-d8	10.646

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
3			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	78
5			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64



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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
Ethanol, 2-(2-b...	10.428	7.4	ng/ul	665439	2	10.646	1788750	20.0