

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN013025\
 Data File : BN036163.D
 Acq On : 31 Jan 2025 11:22
 Operator : RC/JU
 Sample : Q1197-10
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E2985

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_N\Methods\SFAM-EPA-BN013025.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN036163.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.257	8	12	20	rVB	30142	42440	1.34%	0.132%
2	3.675	79	83	101	rBV	129758	215967	6.81%	0.669%
3	3.893	115	120	125	rBV7	2500	4504	0.14%	0.014%
4	3.993	133	137	143	rVB2	7614	11581	0.37%	0.036%
5	4.093	150	154	160	rBV5	8775	13574	0.43%	0.042%
6	4.369	195	201	204	rBV4	2179	3298	0.10%	0.010%
7	4.446	211	214	221	rBV2	1899	3257	0.10%	0.010%
8	4.675	247	253	258	rVB2	6489	10455	0.33%	0.032%
9	4.804	273	275	281	rVB5	2458	4163	0.13%	0.013%
10	4.916	287	294	300	rVB2	7724	13414	0.42%	0.042%
11	5.887	457	459	464	rVB3	2352	3290	0.10%	0.010%
12	6.187	501	510	514	rBV4	7319	11381	0.36%	0.035%
13	6.804	612	615	619	rBV3	3280	4162	0.13%	0.013%
14	6.981	634	645	657	rBV	210042	358078	11.29%	1.110%
15	7.122	661	669	681	rVB	463083	705337	22.24%	2.187%
16	7.334	698	705	714	rBV	581537	893189	28.17%	2.769%
17	7.792	776	783	790	rVB	508501	776797	24.50%	2.408%
18	7.869	790	796	802	rBV	14080	23342	0.74%	0.072%
19	8.516	899	906	914	rBV	529705	853568	26.92%	2.646%
20	8.945	972	979	989	rBV	618573	991167	31.26%	3.073%
21	9.028	989	993	998	rVB3	2941	4305	0.14%	0.013%
22	9.087	998	1003	1008	rBV3	9446	15097	0.48%	0.047%
23	9.392	1051	1055	1061	rBV3	4078	8916	0.28%	0.028%
24	9.510	1072	1075	1080	rVB4	4918	8205	0.26%	0.025%
25	9.669	1094	1102	1115	rBV	561435	904828	28.53%	2.805%
26	9.922	1139	1145	1156	rBV9	6319	20292	0.64%	0.063%
27	10.222	1189	1196	1209	rBV	765564	1247002	39.32%	3.866%
28	10.369	1215	1221	1234	rBV	111835	207662	6.55%	0.644%
29	10.586	1250	1258	1266	rVB	661848	1071907	33.80%	3.323%
30	10.722	1273	1281	1293	rBV	593465	988852	31.18%	3.065%
31	11.133	1343	1351	1358	rBV3	6643	14091	0.44%	0.044%
32	11.239	1365	1369	1375	rVB4	3005	4054	0.13%	0.013%
33	11.392	1387	1395	1400	rBV5	7834	13812	0.44%	0.043%
34	12.086	1508	1513	1518	rBV3	2364	3488	0.11%	0.011%
35	12.175	1523	1528	1533	rVV	9399	14023	0.44%	0.043%
36	12.239	1535	1539	1552	rVB4	3944	10773	0.34%	0.033%

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN013025.MA.M
 Title : SVOA CALIBRATION

37	12.722	1615	1621	1625	rBV5	3662	6434	0.20%	0.020%
38	13.351	1722	1728	1733	rBV4	3806	6677	0.21%	0.021%
39	13.845	1805	1812	1821	rBV	1378428	1853342	58.45%	5.745%
40	13.939	1824	1828	1832	rVV2	10585	15487	0.49%	0.048%
41	14.127	1850	1860	1869	rBV	1360064	1968972	62.09%	6.104%
42	14.433	1906	1912	1919	rVV	995366	1435081	45.26%	4.449%
43	14.516	1922	1926	1930	rVB3	2400	3486	0.11%	0.011%
44	14.669	1945	1952	1960	rBV	215466	333889	10.53%	1.035%
45	14.863	1976	1985	1988	rBV5	3974	8308	0.26%	0.026%
46	15.257	2047	2052	2055	rBV4	3548	4759	0.15%	0.015%
47	15.427	2074	2081	2088	rBV	1833358	2527218	79.70%	7.834%
48	15.539	2093	2100	2111	rBV	817587	1078447	34.01%	3.343%
49	15.669	2118	2122	2126	rBV2	7607	9947	0.31%	0.031%
50	15.792	2134	2143	2148	rBV	106099	139548	4.40%	0.433%
51	15.869	2155	2156	2163	rVB3	3766	4524	0.14%	0.014%
52	15.998	2174	2178	2182	rVB3	2293	3654	0.12%	0.011%
53	16.110	2194	2197	2203	rBV5	2869	4245	0.13%	0.013%
54	16.357	2235	2239	2243	rVB4	10954	13075	0.41%	0.041%
55	16.592	2274	2279	2283	rBV3	4521	6259	0.20%	0.019%
56	16.680	2291	2294	2301	rVB5	4488	6151	0.19%	0.019%
57	16.786	2306	2312	2315	rVB3	2388	3488	0.11%	0.011%
58	16.963	2340	2342	2346	rVB2	4858	4809	0.15%	0.015%
59	17.051	2351	2357	2359	rBV5	3187	4554	0.14%	0.014%
60	17.086	2359	2363	2366	rVB4	4279	4960	0.16%	0.015%
61	17.180	2371	2379	2387	rBV2	1153696	1625566	51.26%	5.039%
62	17.280	2389	2396	2407	rBV2	1956185	2867278	90.42%	8.888%
63	17.851	2490	2493	2497	rVB5	6290	7819	0.25%	0.024%
64	17.951	2504	2510	2517	rBV9	5726	15610	0.49%	0.048%
65	18.080	2527	2532	2541	rBV	73218	99271	3.13%	0.308%
66	18.410	2580	2588	2593	rBV	10847	27760	0.88%	0.086%
67	18.504	2602	2604	2608	rBV3	2467	3195	0.10%	0.010%
68	18.562	2608	2614	2616	rVV6	6063	9620	0.30%	0.030%
69	18.933	2670	2677	2681	rBV6	8563	16982	0.54%	0.053%
70	19.139	2708	2712	2716	rVB2	26535	34214	1.08%	0.106%
71	19.210	2719	2724	2728	rBV2	40662	66171	2.09%	0.205%
72	19.251	2728	2731	2739	rVB8	17093	37202	1.17%	0.115%
73	19.362	2746	2750	2754	rBV6	24642	32113	1.01%	0.100%
74	19.515	2773	2776	2780	rBV4	12681	14212	0.45%	0.044%
75	19.574	2780	2786	2793	rBV	2398628	3171031	100.00%	9.830%
76	21.115	3044	3048	3053	rVB5	38867	56043	1.77%	0.174%

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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_N\Methods\SFAM-EPA-BN013025.MA.M
Title : SVOA CALIBRATION

77	21.427	3095	3101	3110	rBV	1015642	1518060	47.87%	4.706%
78	24.245	3571	3580	3592	rBV	989586	2325284	73.33%	7.208%
79	24.456	3608	3616	3634	rVB2	511812	1403315	44.25%	4.350%

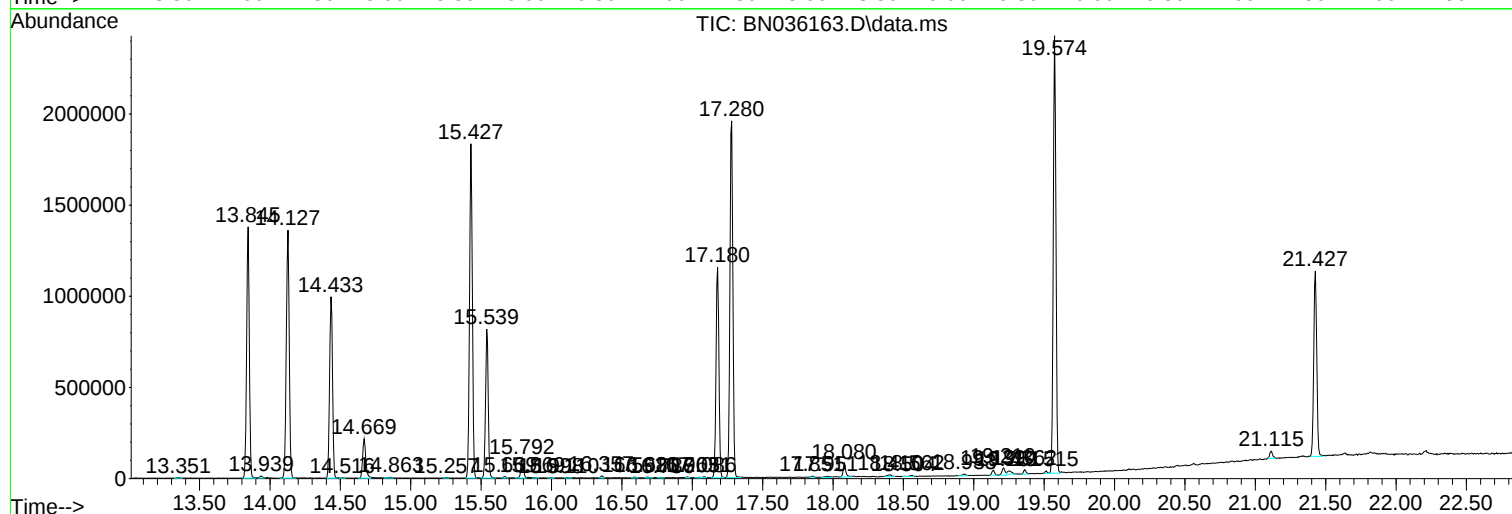
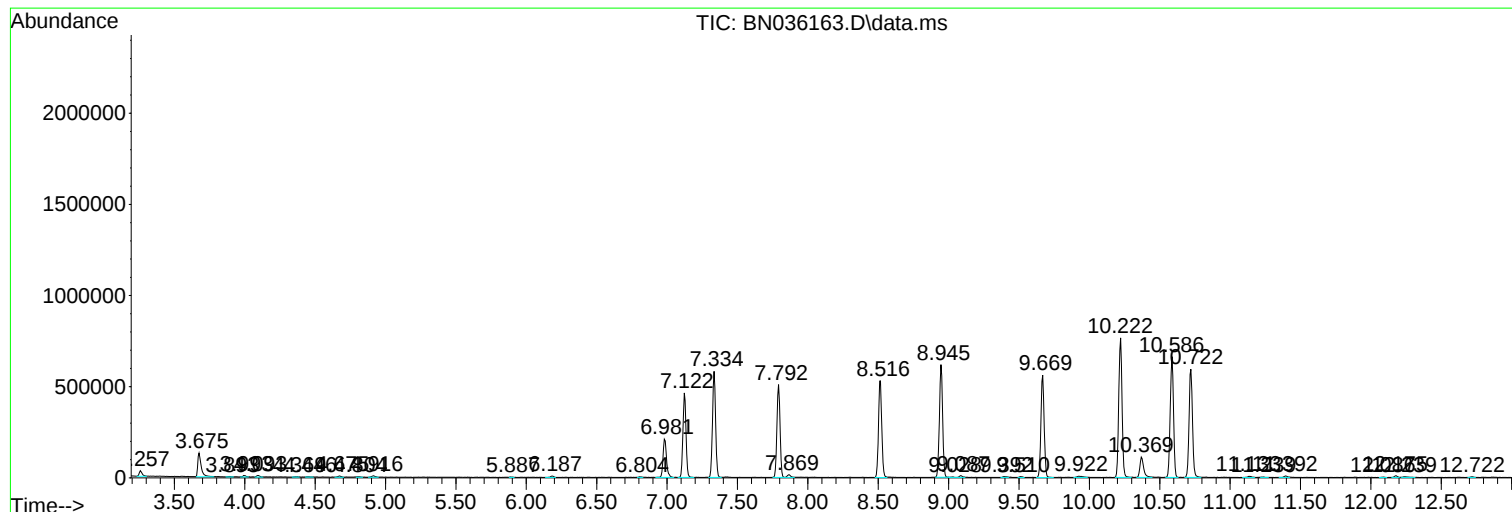
Sum of corrected areas: 32258331

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN013025.MA.M
Quant Title : SVOA CALIBRATION

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



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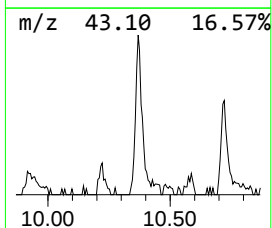
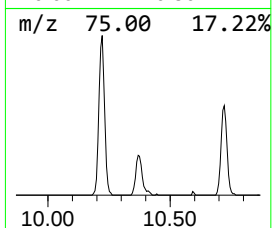
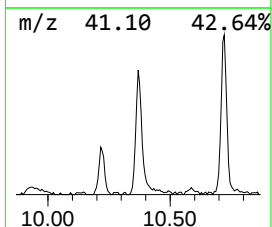
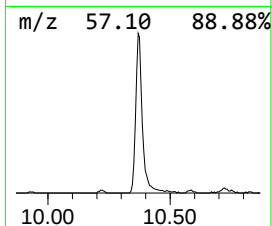
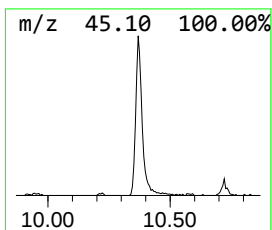
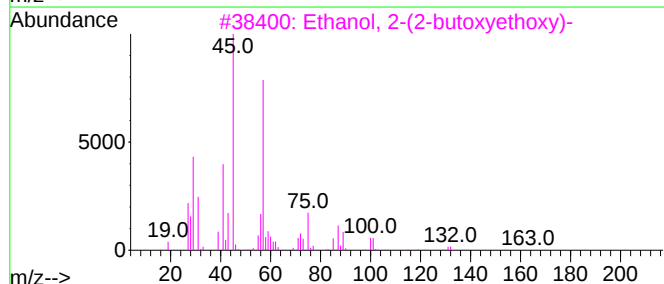
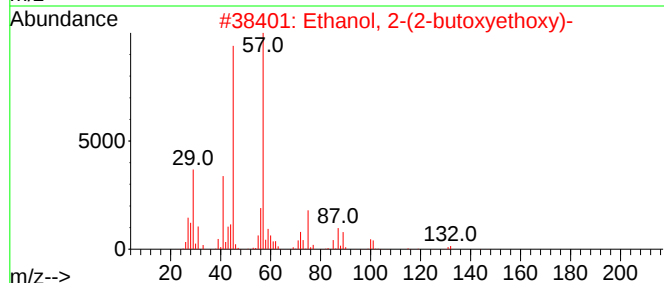
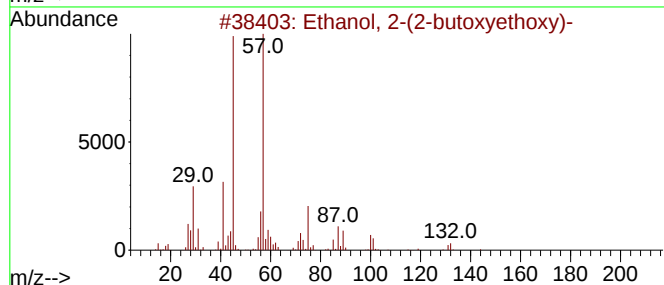
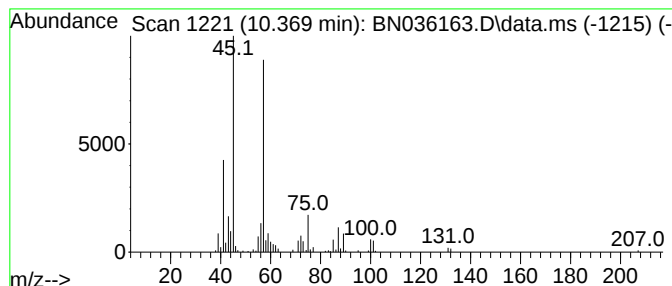
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN013025.MA.M
Quant Title : SVOA CALIBRATION

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.
10.369	3.87 ng/ul	207662	Naphthalene-d8	10.586

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	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	78



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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...	10.369	3.9	ng/u1	207662	2	10.586	1071910	20.0