

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101822\
 Data File : BN022152.D
 Acq On : 17 Oct 2022 15:24
 Operator : CG/JU
 Sample : N5070-14
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BS7

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-BN101222.M
 Title : SVOA CALIBRATION

Signal : TIC: BN022152.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.258	42	46	51	rBV	22051	32514	0.74%	0.084%
2	3.422	71	74	79	rVB3	4377	6217	0.14%	0.016%
3	3.546	92	95	113	rBV	111770	187116	4.27%	0.485%
4	3.699	119	121	135	rBV	114529	204584	4.67%	0.530%
5	3.934	155	161	169	rVB2	13206	26284	0.60%	0.068%
6	4.034	173	178	183	rBV	8490	13259	0.30%	0.034%
7	4.422	242	244	249	rVB4	2865	4429	0.10%	0.011%
8	7.110	695	701	709	rBV	127536	205385	4.69%	0.532%
9	7.281	723	730	744	rBV	506533	802690	18.34%	2.081%
10	7.475	756	763	774	rBV	458309	740602	16.92%	1.920%
11	7.557	774	777	786	rVB7	3644	6285	0.14%	0.016%
12	7.952	836	844	853	rBV	537350	871280	19.90%	2.259%
13	8.675	960	967	981	rVB	377133	615872	14.07%	1.597%
14	9.134	1038	1045	1058	rBV	623749	1022322	23.35%	2.650%
15	9.863	1161	1169	1178	rBV	390429	654591	14.95%	1.697%
16	10.404	1253	1261	1280	rBV	670401	1157193	26.43%	3.000%
17	10.787	1318	1326	1334	rBV	817396	1414463	32.31%	3.667%
18	10.928	1339	1350	1366	rBV	617346	1048529	23.95%	2.718%
19	12.034	1534	1538	1545	rVB4	3269	6634	0.15%	0.017%
20	12.381	1592	1597	1603	rBV	12317	20488	0.47%	0.053%
21	13.592	1796	1803	1810	rBV5	3236	6764	0.15%	0.018%
22	14.039	1871	1879	1888	rBV	1421591	1934658	44.19%	5.016%
23	14.322	1920	1927	1940	rBV2	1617054	2377038	54.30%	6.162%
24	14.586	1968	1972	1973	rBV3	3755	4585	0.10%	0.012%
25	14.628	1973	1979	1986	rBV	1346239	1961452	44.80%	5.085%
26	14.834	2008	2014	2028	rBV	89612	167749	3.83%	0.435%
27	15.051	2046	2051	2058	rVB4	20161	34690	0.79%	0.090%
28	15.134	2060	2065	2070	rBV7	4339	8096	0.18%	0.021%
29	15.410	2109	2112	2118	rVB7	4355	6887	0.16%	0.018%
30	15.528	2126	2132	2136	rBV6	2766	5367	0.12%	0.014%
31	15.622	2141	2148	2158	rVV	2175644	3096737	70.74%	8.028%
32	15.745	2163	2169	2181	rVV	435354	592048	13.52%	1.535%
33	15.869	2184	2190	2195	rVV5	17878	33629	0.77%	0.087%
34	15.969	2204	2207	2212	rBV7	3570	4516	0.10%	0.012%
35	16.086	2223	2227	2232	rVB7	3333	6338	0.14%	0.016%
36	16.275	2254	2259	2261	rBV5	3697	6866	0.16%	0.018%

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37	16.351	2267	2272	2277	rBV	22837	36497	0.83%	0.095%
38	16.675	2323	2327	2331	rBV6	5022	8295	0.19%	0.022%
39	17.180	2407	2413	2417	rBV3	23190	38098	0.87%	0.099%
40	17.386	2441	2448	2454	rBV	1663927	2410061	55.05%	6.248%
41	17.486	2458	2465	2476	rVV2	2700290	3970830	90.70%	10.294%
42	18.269	2595	2598	2603	rBV3	17230	22495	0.51%	0.058%
43	19.345	2777	2781	2787	rVB	37105	54893	1.25%	0.142%
44	19.416	2789	2793	2799	rVB	40904	54343	1.24%	0.141%
45	19.698	2839	2841	2846	rBV5	7673	10238	0.23%	0.027%
46	19.786	2849	2856	2864	rBV2	3168039	4377838	100.00%	11.349%
47	21.580	3156	3161	3169	rBV	1754194	2421173	55.31%	6.277%
48	23.892	3546	3554	3572	rVB2	1807723	3782888	86.41%	9.807%
49	24.051	3574	3581	3592	rVB	1036537	2097647	47.92%	5.438%

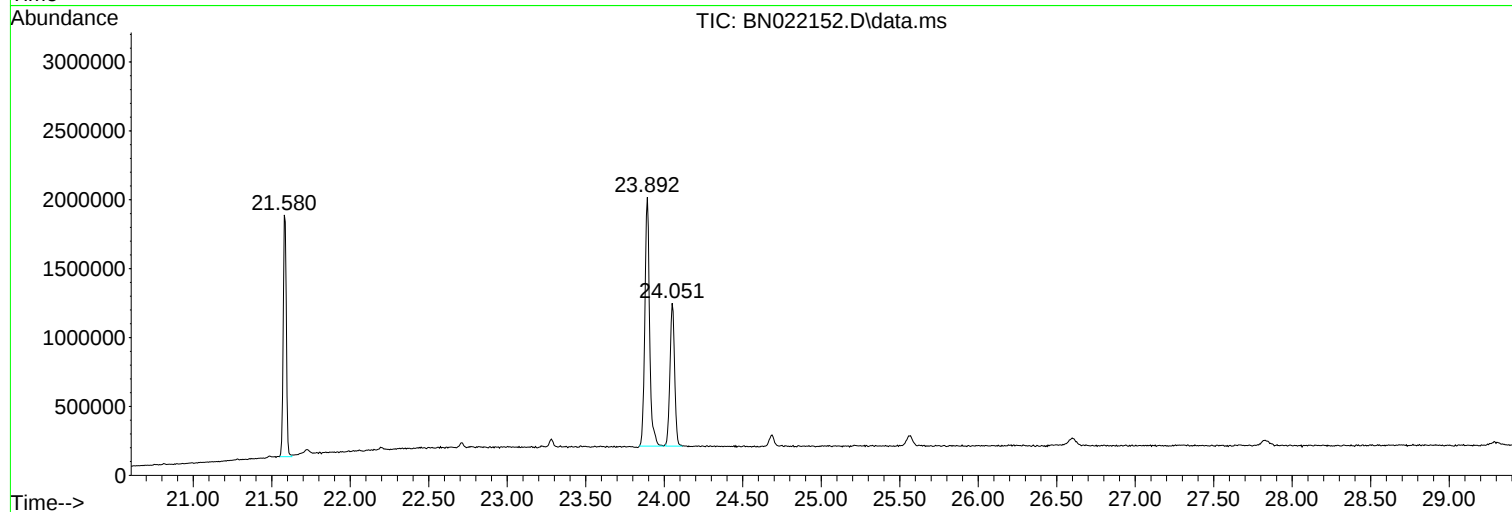
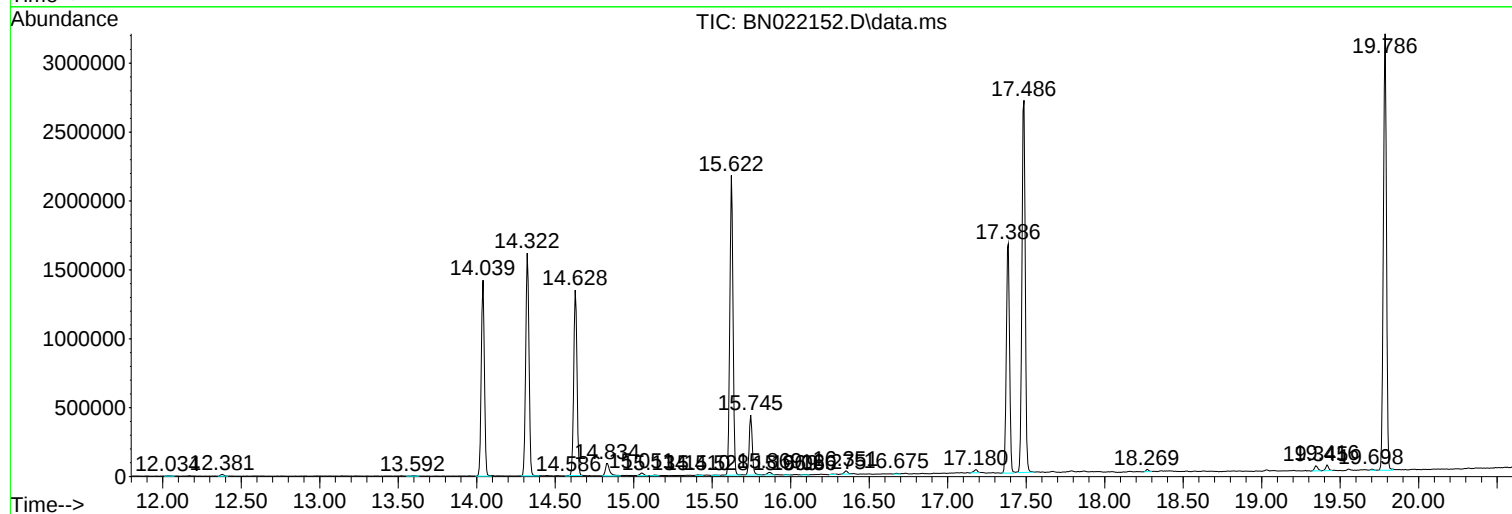
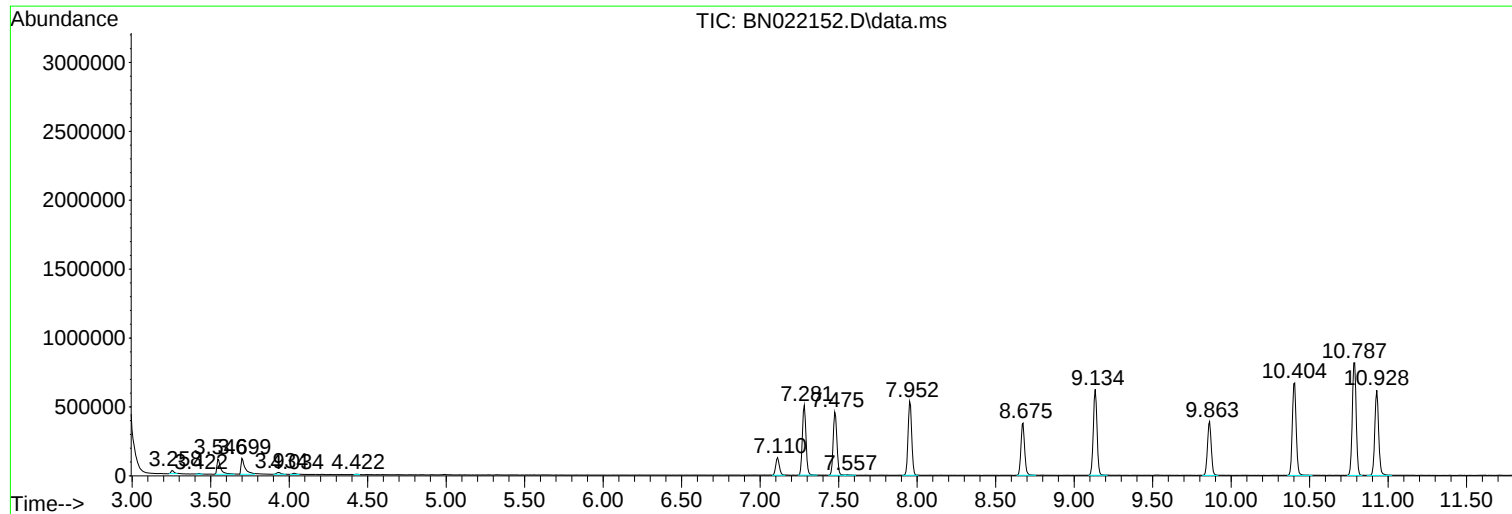
Sum of corrected areas: 38573453

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101222.M
Quant Title : SVOA CALIBRATION

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



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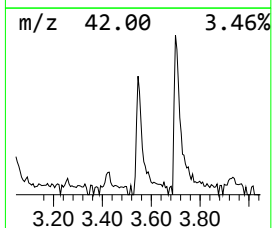
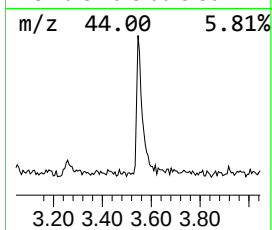
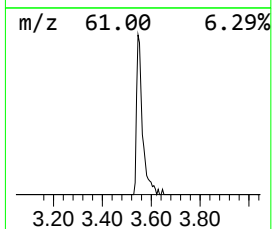
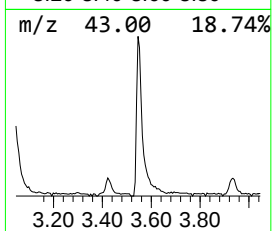
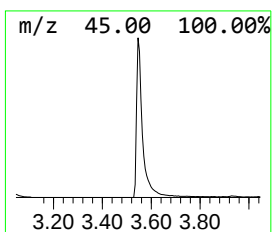
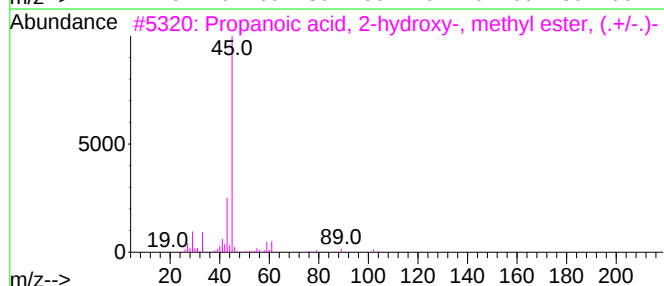
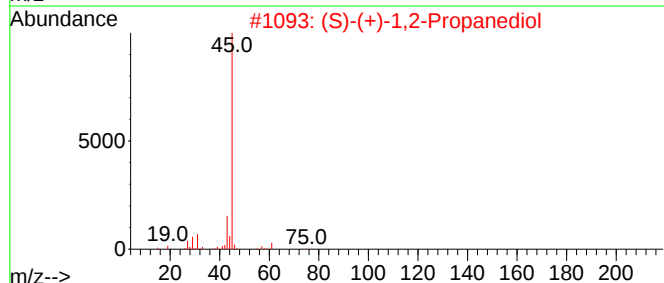
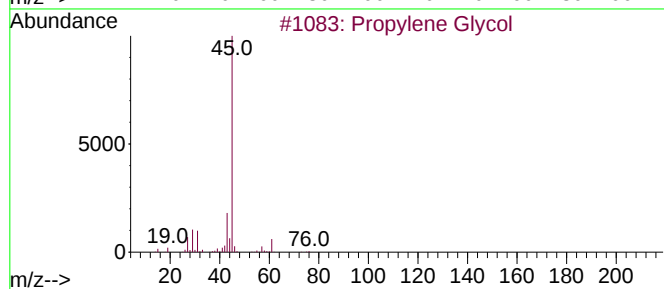
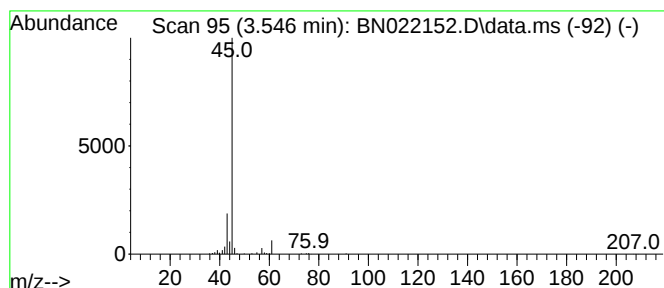
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.546	4.30 ng/ul	187116	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	90
2			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	56
3			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
4			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	37
5			Isopropyl Alcohol	60	C3H8O	000067-63-0	9



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propylene Glycol	3.546	4.3	ng/u1	187116	1	7.952	871280	20.0