

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
 Data File : BP007781.D
 Acq On : 29 Oct 2021 12:55
 Operator : CG/JU
 Sample : M4281-18DL 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFY88DL

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-BP102521.M
 Title : SVOA CALIBRATION

Signal : TIC: BP007781.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.340	6	9	13	rBV5	15090	25469	0.27%	0.053%
2	3.405	16	20	30	rVB	201314	286238	3.07%	0.597%
3	3.675	55	66	70	rBV	4086144	9313505	100.00%	19.429%
4	3.799	85	87	95	rVV	143095	191275	2.05%	0.399%
5	3.893	99	103	111	rVB2	35146	60067	0.64%	0.125%
6	4.028	123	126	129	rVB	13789	16999	0.18%	0.035%
7	4.116	138	141	147	rVB3	15474	21160	0.23%	0.044%
8	4.205	154	156	162	rVB7	8009	13861	0.15%	0.029%
9	4.887	268	272	277	rBV2	15599	24392	0.26%	0.051%
10	6.322	512	516	520	rVB7	7035	9722	0.10%	0.020%
11	6.928	614	619	626	rBV8	9104	22044	0.24%	0.046%
12	7.093	642	647	657	rVB	109591	184361	1.98%	0.385%
13	7.257	669	675	689	rVB	305123	486647	5.23%	1.015%
14	7.463	703	710	718	rBV	336720	529738	5.69%	1.105%
15	7.852	769	776	780	rBV5	10542	19572	0.21%	0.041%
16	7.934	783	790	798	rVV	879270	1429099	15.34%	2.981%
17	8.640	903	910	922	rBV	289845	467378	5.02%	0.975%
18	9.087	979	986	995	rVB	352209	575816	6.18%	1.201%
19	9.557	1059	1066	1074	rVB9	12791	29594	0.32%	0.062%
20	9.810	1103	1109	1117	rBV	267626	447706	4.81%	0.934%
21	10.357	1195	1202	1213	rBV	496812	842141	9.04%	1.757%
22	10.516	1223	1229	1237	rBV2	92500	156219	1.68%	0.326%
23	10.734	1259	1266	1279	rVB	1157899	1972797	21.18%	4.116%
24	10.869	1281	1289	1298	rBV	348564	615026	6.60%	1.283%
25	12.263	1517	1526	1531	rBV2	37073	58875	0.63%	0.123%
26	12.322	1531	1536	1542	rVB4	9368	16552	0.18%	0.035%
27	12.940	1638	1641	1646	rVB7	8253	12056	0.13%	0.025%
28	13.192	1679	1684	1689	rVB6	7633	10981	0.12%	0.023%
29	13.475	1728	1732	1735	rBV4	8031	12375	0.13%	0.026%
30	13.875	1795	1800	1804	rBV8	5929	12961	0.14%	0.027%
31	13.981	1810	1818	1824	rBV	956400	1314084	14.11%	2.741%
32	14.263	1855	1866	1878	rBV	966564	1436718	15.43%	2.997%
33	14.487	1897	1904	1909	rBV	23458	57324	0.62%	0.120%
34	14.569	1911	1918	1938	rVB	1792445	2791394	29.97%	5.823%
35	14.769	1946	1952	1961	rBV	85252	147952	1.59%	0.309%
36	14.916	1973	1977	1986	rVB9	14089	28823	0.31%	0.060%

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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-BP102521.M
 Title : SVOA CALIBRATION

37	14.998	1986	1991	1992	rBV5	9758	13186	0.14%	0.028%
38	15.022	1992	1995	2004	rVV	17209	29050	0.31%	0.061%
39	15.110	2004	2010	2018	rVV4	19607	46176	0.50%	0.096%
40	15.234	2026	2031	2035	rBV7	7306	11253	0.12%	0.023%
41	15.386	2052	2057	2061	rBV2	27704	43868	0.47%	0.092%
42	15.434	2061	2065	2071	rVB2	19402	30258	0.32%	0.063%
43	15.569	2081	2088	2095	rBV	1338384	1989777	21.36%	4.151%
44	15.681	2102	2107	2114	rBV	227132	320729	3.44%	0.669%
45	15.810	2123	2129	2135	rVB2	19349	29705	0.32%	0.062%
46	15.939	2144	2151	2153	rBV8	6672	11720	0.13%	0.024%
47	16.010	2160	2163	2169	rVB8	8553	15201	0.16%	0.032%
48	16.145	2179	2186	2195	rBV	15345	37666	0.40%	0.079%
49	16.251	2200	2204	2209	rVV8	9293	17906	0.19%	0.037%
50	16.298	2209	2212	2215	rVV5	7347	11656	0.13%	0.024%
51	16.363	2219	2223	2228	rVB7	7023	12664	0.14%	0.026%
52	16.639	2266	2270	2277	rVV8	11227	20221	0.22%	0.042%
53	16.980	2324	2328	2337	rBV	33300	65916	0.71%	0.138%
54	17.110	2347	2350	2353	rBV2	10255	14636	0.16%	0.031%
55	17.139	2353	2355	2361	rVB5	15596	22494	0.24%	0.047%
56	17.210	2363	2367	2369	rBV5	5973	9640	0.10%	0.020%
57	17.328	2381	2387	2393	rBV	2391793	3436350	36.90%	7.169%
58	17.428	2398	2404	2415	rVB	1598785	2256923	24.23%	4.708%
59	17.645	2435	2441	2444	rBV2	12713	20287	0.22%	0.042%
60	18.010	2499	2503	2507	rVB2	35048	44115	0.47%	0.092%
61	18.222	2534	2539	2548	rBV	151407	230407	2.47%	0.481%
62	18.686	2615	2618	2619	rBV3	11122	12707	0.14%	0.027%
63	18.951	2659	2663	2666	rBV	32065	52410	0.56%	0.109%
64	19.310	2721	2724	2725	rBV2	23067	24187	0.26%	0.050%
65	19.339	2725	2729	2734	rVB	133982	163644	1.76%	0.341%
66	19.457	2747	2749	2753	rVB4	28158	28222	0.30%	0.059%
67	19.604	2772	2774	2778	rVB3	29190	30209	0.32%	0.063%
68	19.663	2778	2784	2795	rBV	2204158	3074747	33.01%	6.414%
69	21.133	3031	3034	3040	rBV5	168272	223755	2.40%	0.467%
70	21.421	3077	3083	3088	rBV	3375427	4509315	48.42%	9.407%
71	23.710	3465	3472	3480	rVB2	1537646	2859355	30.70%	5.965%
72	23.868	3492	3499	3507	rVB2	2295546	4573757	49.11%	9.542%

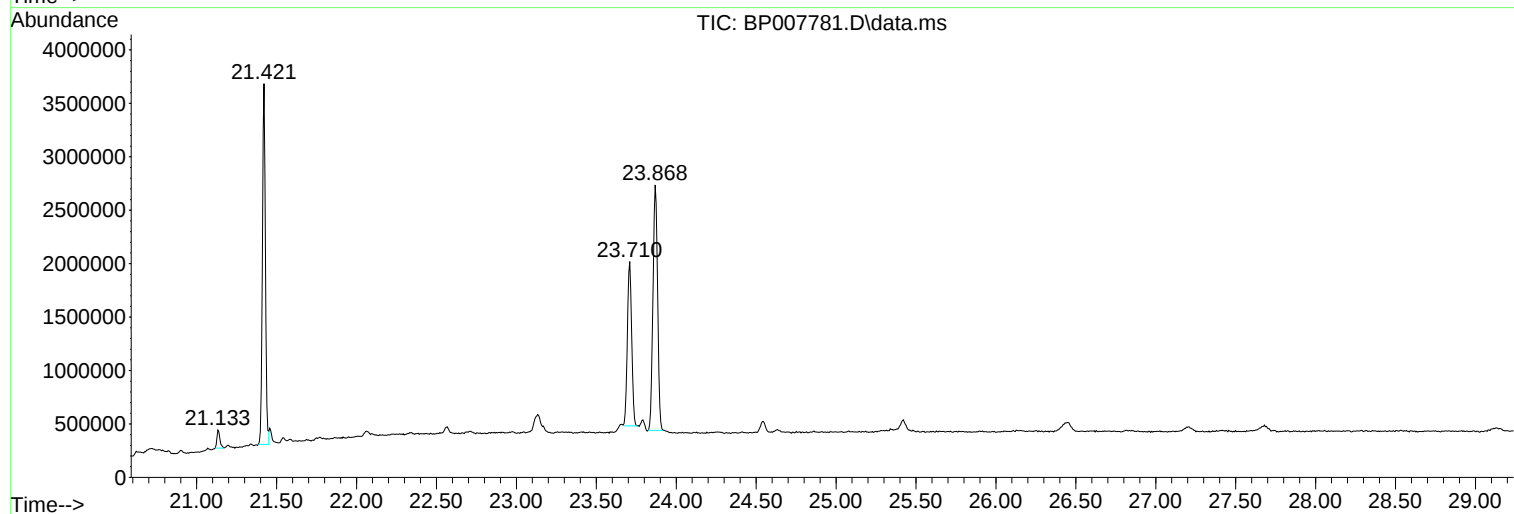
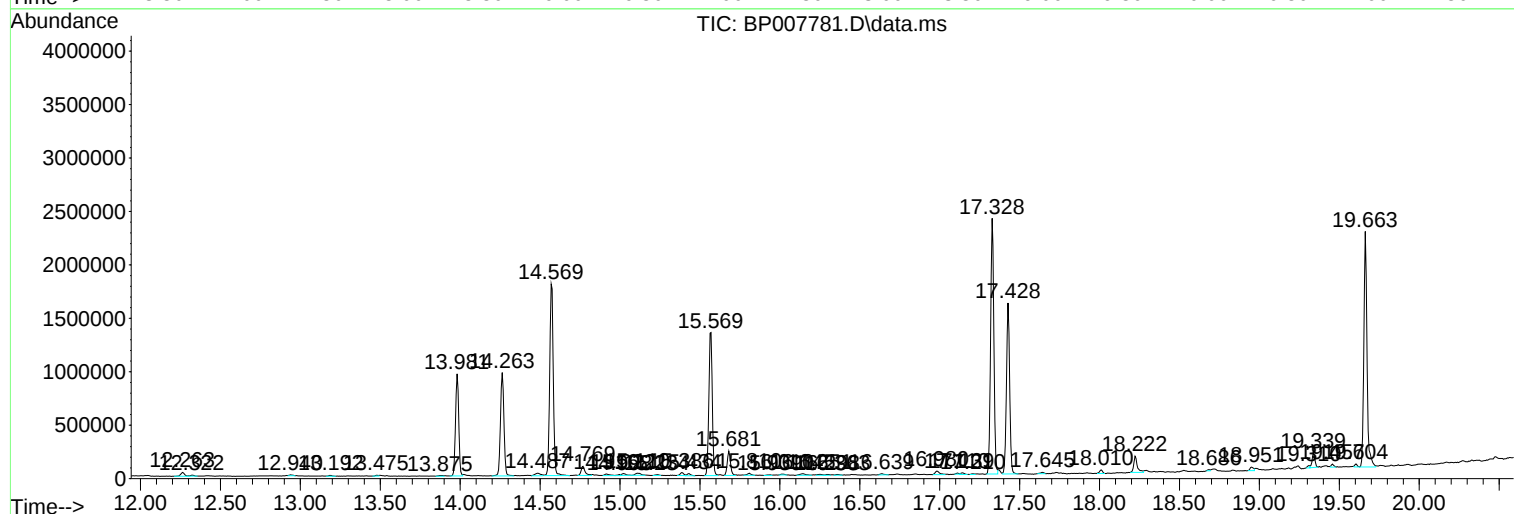
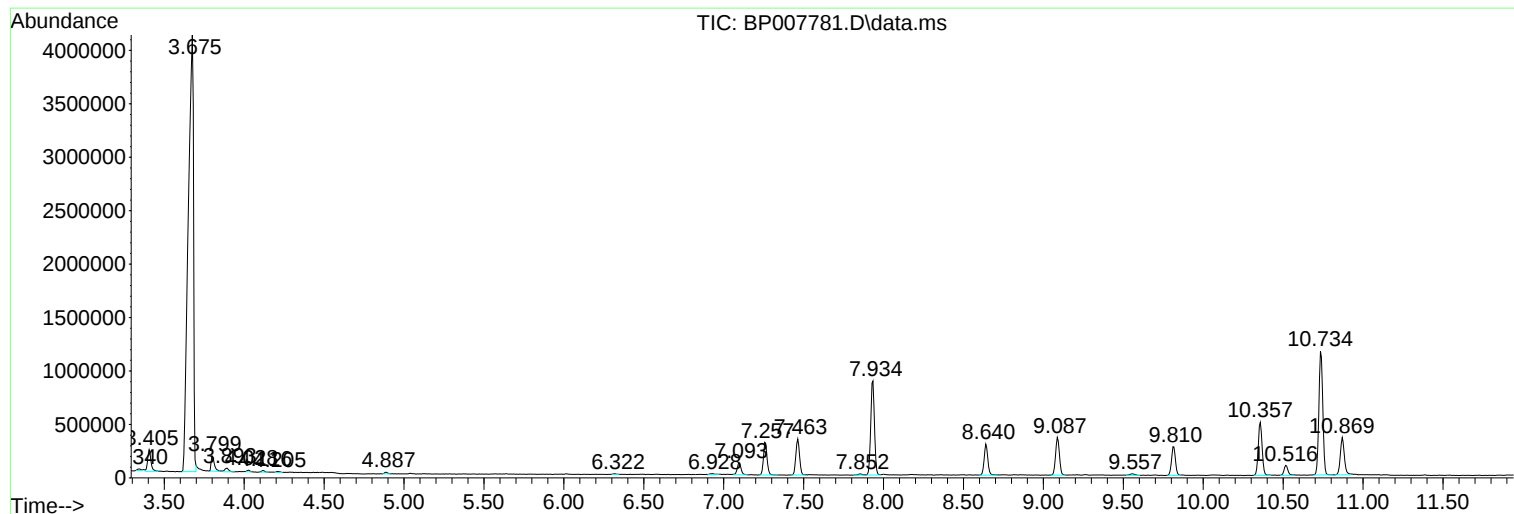
Sum of corrected areas: 47935033

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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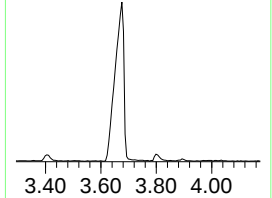
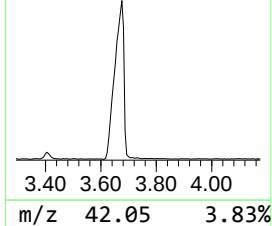
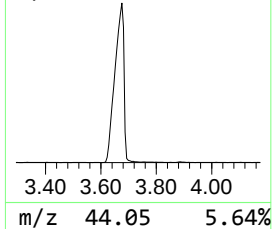
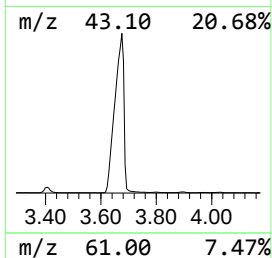
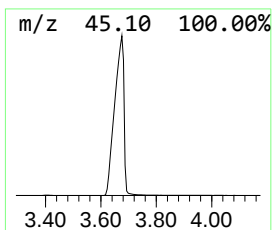
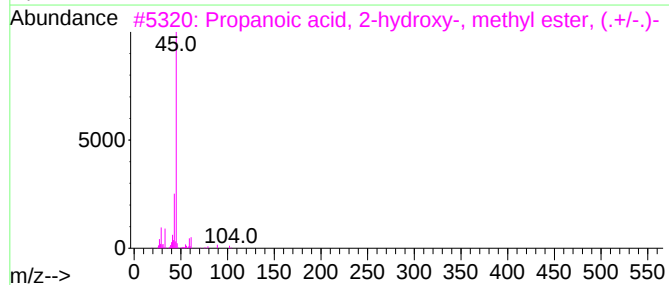
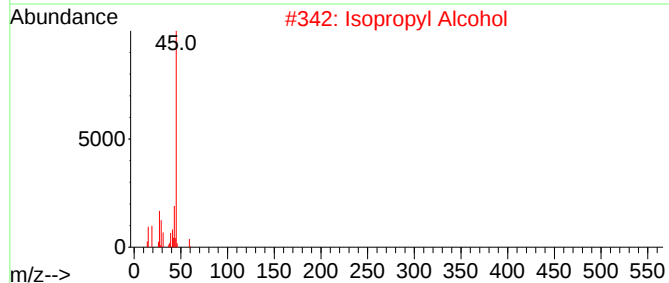
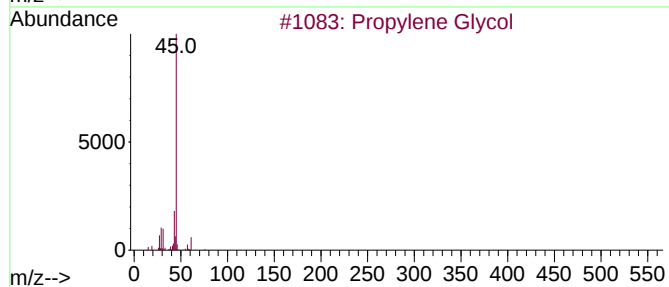
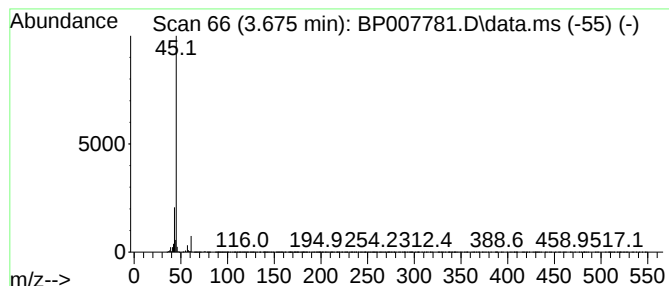
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

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.
3.675	130.34 ng/ul	9313510	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	90
2			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
3			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
4			Isopropyl Alcohol	60	C3H8O	000067-63-0	9
5			Propane, 2-ethoxy-	88	C5H12O	000625-54-7	9



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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propylene Glycol	3.675	130.3	ng/ul	9313510	1	7.934	1429100	20.0