

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
 Data File : BP007611.D
 Acq On : 23 Oct 2021 03:41
 Operator : CG/JU
 Sample : M4281-19
 Misc :
 ALS Vial : 71 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFY97

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-BP101421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP007611.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	11	15	20	rBV	40827	56683	0.90%	0.103%
2	3.540	39	43	49	rVB8	5846	9811	0.15%	0.018%
3	3.646	57	61	70	rBV	99939	161997	2.56%	0.294%
4	3.799	84	87	96	rBV	92991	147148	2.32%	0.267%
5	4.034	123	127	132	rVB	31518	44943	0.71%	0.082%
6	4.122	139	142	148	rVB	14344	18105	0.29%	0.033%
7	4.222	154	159	164	rVB5	13190	20277	0.32%	0.037%
8	5.052	295	300	306	rVB3	11778	17819	0.28%	0.032%
9	5.605	389	394	399	rBV8	4456	6827	0.11%	0.012%
10	6.328	511	517	522	rBV2	13101	19618	0.31%	0.036%
11	7.105	641	649	658	rVB	225793	377545	5.96%	0.685%
12	7.269	671	677	689	rVB	727948	1149171	18.15%	2.085%
13	7.475	703	712	721	rBV	799965	1264326	19.97%	2.294%
14	7.946	784	792	800	rBV	729040	1162506	18.36%	2.109%
15	8.016	800	804	809	rVB7	6392	10043	0.16%	0.018%
16	8.199	829	835	840	rBV9	5118	9560	0.15%	0.017%
17	8.505	882	887	890	rBV7	4426	7157	0.11%	0.013%
18	8.652	903	912	924	rBV	630296	1015356	16.04%	1.842%
19	9.099	980	988	997	rBV	892794	1446735	22.85%	2.625%
20	9.563	1061	1067	1074	rVB	24549	41868	0.66%	0.076%
21	9.828	1102	1112	1121	rBV	673839	1146488	18.11%	2.080%
22	10.369	1196	1204	1218	rBV	1092415	1845002	29.15%	3.348%
23	10.528	1226	1231	1237	rBV	46720	74807	1.18%	0.136%
24	10.746	1260	1268	1280	rVB	965289	1669402	26.37%	3.029%
25	10.881	1283	1291	1311	rVV	748408	1295873	20.47%	2.351%
26	12.063	1486	1492	1498	rVB3	17257	30031	0.47%	0.054%
27	12.257	1521	1525	1532	rVB10	7077	15215	0.24%	0.028%
28	12.334	1532	1538	1545	rBV2	24320	42783	0.68%	0.078%
29	13.493	1731	1735	1739	rVB6	5227	7766	0.12%	0.014%
30	13.551	1739	1745	1750	rBV8	4980	8103	0.13%	0.015%
31	13.993	1813	1820	1826	rBV	2058339	2878995	45.48%	5.224%
32	14.128	1839	1843	1849	rVB8	7385	13609	0.21%	0.025%
33	14.275	1857	1868	1876	rBV	2304650	3637695	57.46%	6.601%
34	14.504	1902	1907	1912	rVV4	9845	17982	0.28%	0.033%
35	14.581	1913	1920	1938	rVB	1550041	2326025	36.74%	4.221%
36	14.769	1943	1952	1965	rBV	210401	340737	5.38%	0.618%

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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-BP101421.M
 Title : SVOA CALIBRATION

37	14.998	1986	1991	1995	rBV8	10249	16230	0.26%	0.029%
38	15.116	2009	2011	2017	rVB7	5754	8422	0.13%	0.015%
39	15.240	2025	2032	2036	rBV5	20597	34919	0.55%	0.063%
40	15.328	2042	2047	2050	rBV	14376	20957	0.33%	0.038%
41	15.398	2055	2059	2064	rBV4	10190	15189	0.24%	0.028%
42	15.575	2081	2089	2098	rBV	3118268	4449029	70.28%	8.073%
43	15.687	2102	2108	2117	rVV	407759	565701	8.94%	1.026%
44	15.816	2123	2130	2135	rVB	41360	63352	1.00%	0.115%
45	16.145	2181	2186	2191	rBV9	5911	11771	0.19%	0.021%
46	16.251	2203	2204	2209	rVB5	6959	8055	0.13%	0.015%
47	17.004	2326	2332	2337	rBV9	9260	23496	0.37%	0.043%
48	17.116	2346	2351	2354	rBV3	12203	22937	0.36%	0.042%
49	17.334	2380	2388	2395	rBV	1901816	2736190	43.22%	4.965%
50	17.434	2398	2405	2417	rBV2	3570400	5198144	82.11%	9.432%
51	17.528	2418	2421	2425	rVV	20056	25294	0.40%	0.046%
52	17.575	2426	2429	2435	rVB5	12360	17722	0.28%	0.032%
53	17.710	2449	2452	2456	rBV5	6789	8805	0.14%	0.016%
54	17.975	2495	2497	2499	rBV3	11463	11621	0.18%	0.021%
55	18.010	2499	2503	2512	rVB4	54120	76669	1.21%	0.139%
56	18.228	2536	2540	2547	rBV	89935	126869	2.00%	0.230%
57	19.239	2709	2712	2717	rBV2	48953	65385	1.03%	0.119%
58	19.310	2720	2724	2728	rBV	58319	75864	1.20%	0.138%
59	19.457	2746	2749	2754	rBV4	32133	43424	0.69%	0.079%
60	19.663	2778	2784	2792	rBV	4533482	6152485	97.19%	11.164%
61	21.133	3031	3034	3040	rBV3	141612	186460	2.95%	0.338%
62	21.416	3077	3082	3088	rBV	2379144	3214153	50.77%	5.832%
63	23.710	3464	3472	3480	rVB2	3072205	6330390	100.00%	11.486%
64	23.863	3491	3498	3509	rVB2	1579235	3264115	51.56%	5.923%

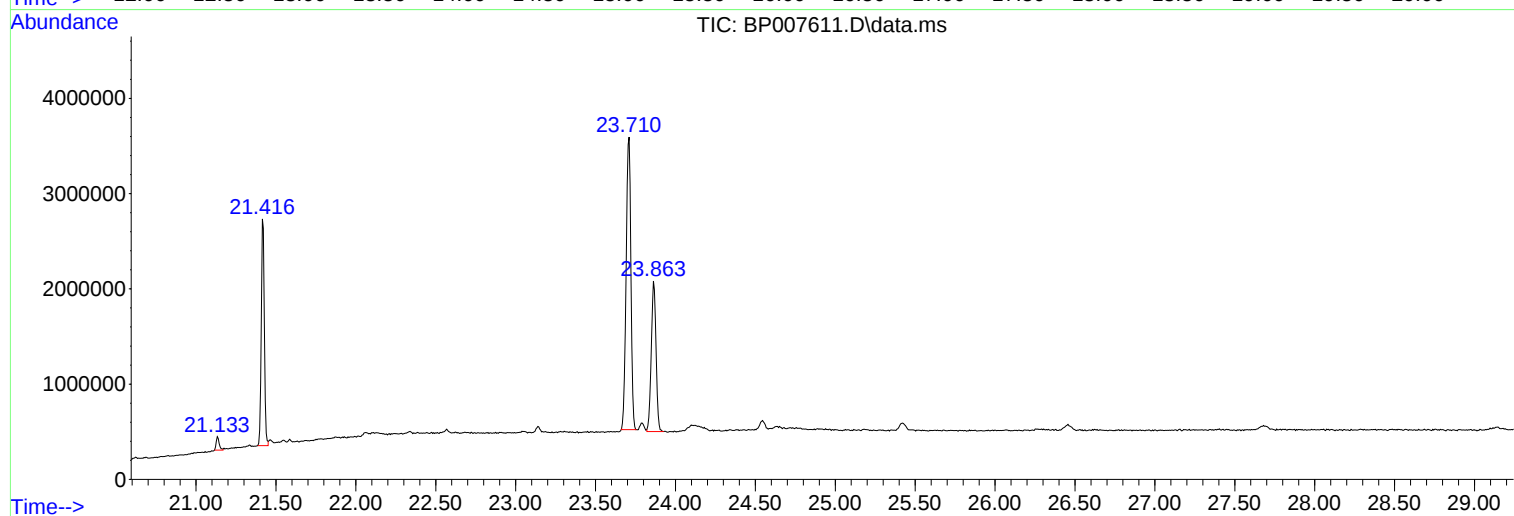
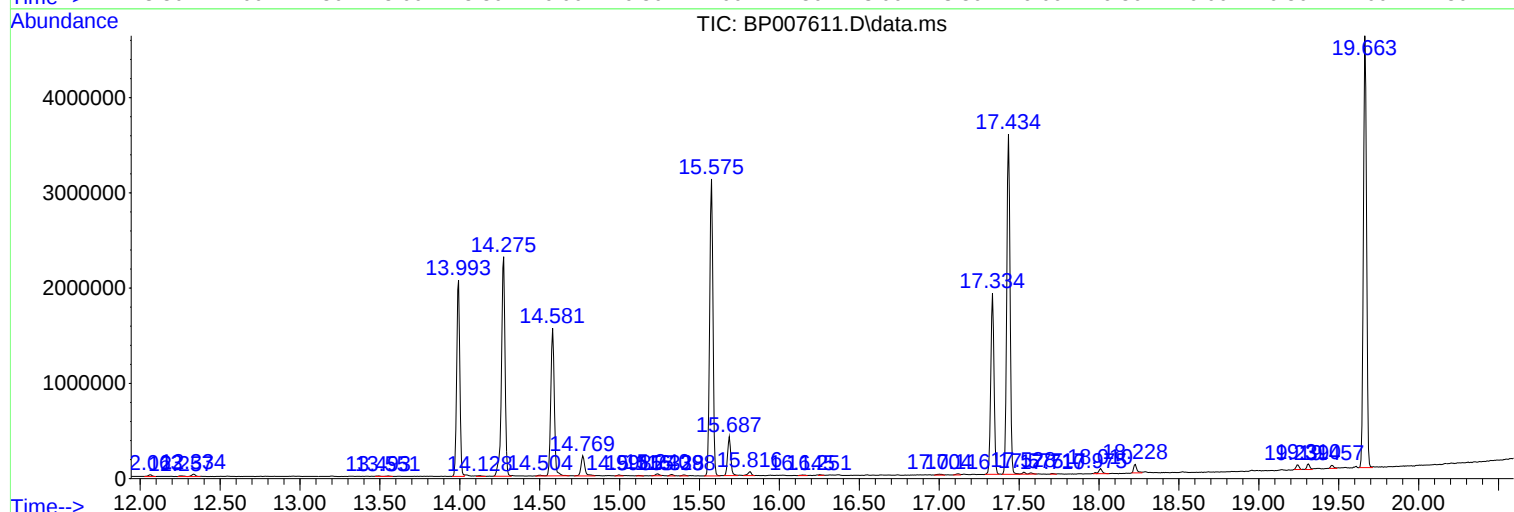
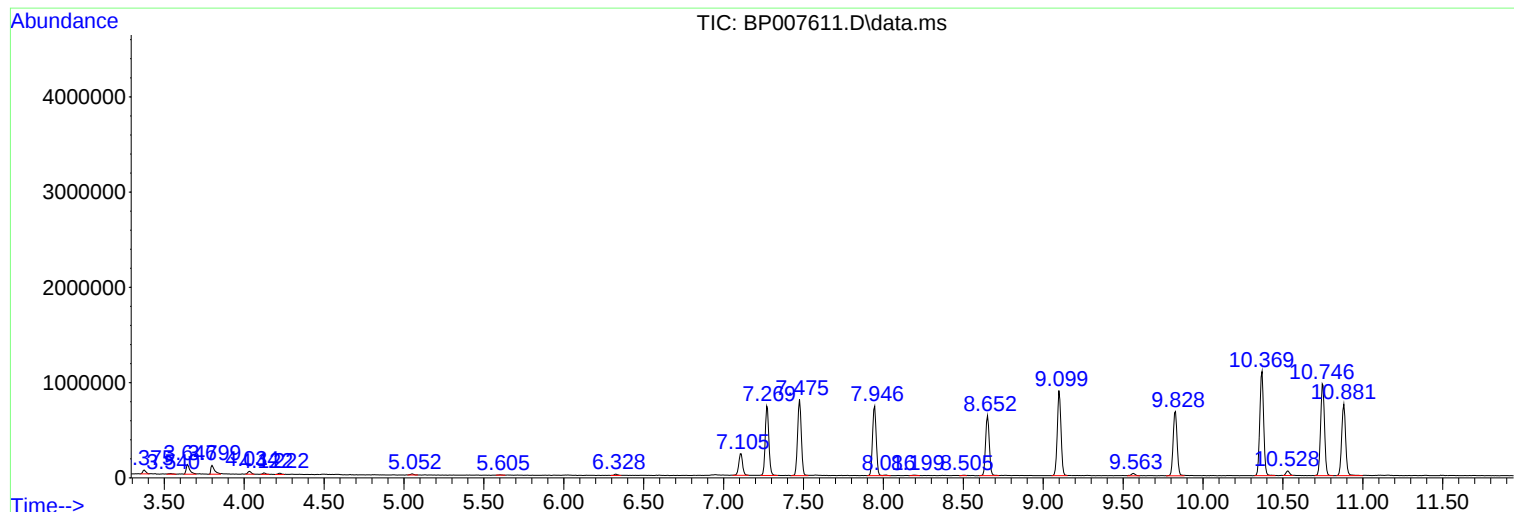
Sum of corrected areas: 55111656

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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



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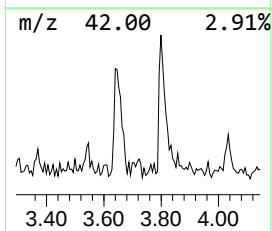
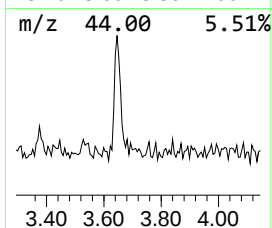
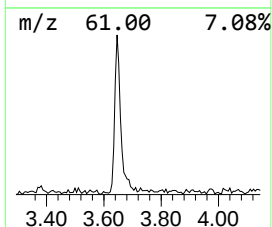
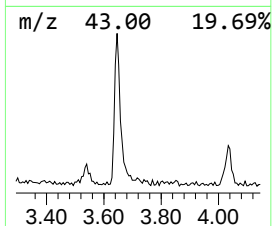
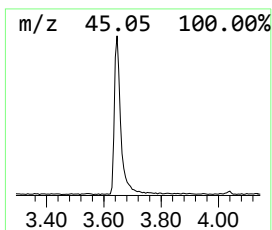
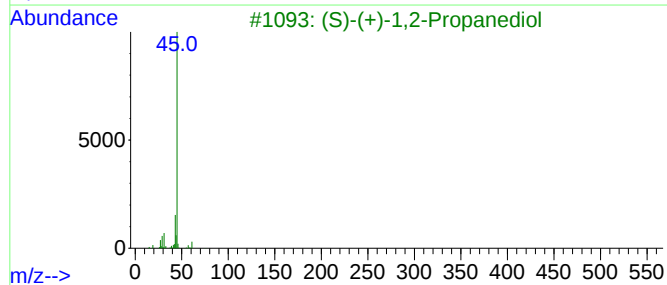
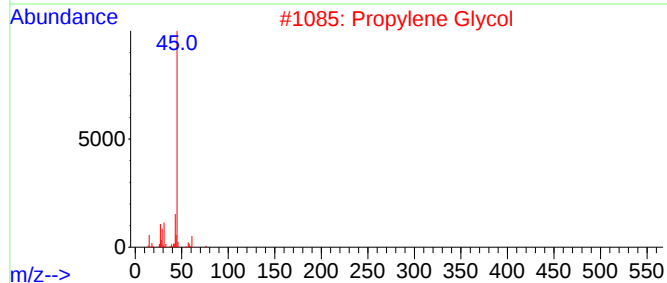
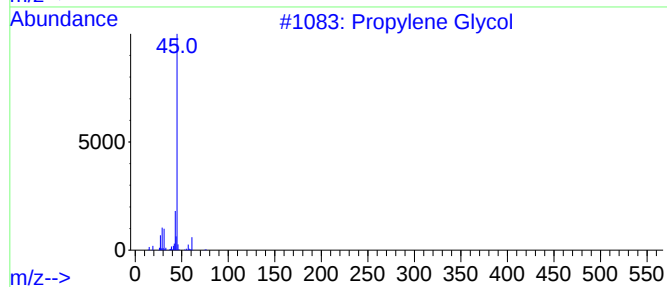
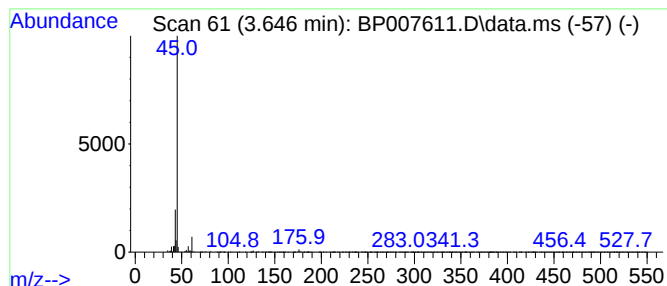
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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.
3.646	2.79 ng/ul	161997	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	72
2			Propylene Glycol	76	C3H8O2	000057-55-6	40
3			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	39
4			Isopropyl Alcohol	60	C3H8O	000067-63-0	9
5			Isopropyl Alcohol	60	C3H8O	000067-63-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propylene Glycol	3.646	2.8	ng/ul	161997	1	7.946	1162510	20.0