

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
 Data File : BP021679.D
 Acq On : 27 Aug 2024 20:28
 Operator : RC/JU
 Sample : P3675-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCYE2

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-BP082324.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021679.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.017	3	5	21	rVB	2136112	2690044	47.01%	3.945%
2	3.275	45	49	58	rVB	43909	59505	1.04%	0.087%
3	3.540	90	94	106	rVB	176190	244879	4.28%	0.359%
4	3.687	115	119	133	rVB	564598	821180	14.35%	1.204%
5	4.016	172	175	181	rVB	24887	34781	0.61%	0.051%
6	4.675	283	287	291	rBV4	10093	16609	0.29%	0.024%
7	4.922	326	329	335	rVB2	10759	15502	0.27%	0.023%
8	5.893	490	494	498	rBV	23156	31171	0.54%	0.046%
9	6.063	519	523	531	rVB10	6595	11230	0.20%	0.016%
10	6.505	592	598	603	rVB5	10328	17560	0.31%	0.026%
11	6.793	640	647	654	rBV2	15962	34946	0.61%	0.051%
12	6.957	666	675	691	rBV	740230	1234391	21.57%	1.810%
13	7.122	696	703	715	rVB	638470	1004469	17.55%	1.473%
14	7.328	730	738	744	rBV	702808	1141726	19.95%	1.674%
15	7.393	745	749	759	rVB	115398	188230	3.29%	0.276%
16	7.699	798	801	809	rBV10	3637	8511	0.15%	0.012%
17	7.793	810	817	824	rVV	1228835	1942657	33.95%	2.849%
18	7.857	824	828	833	rVB2	18887	30862	0.54%	0.045%
19	8.157	877	879	885	rVB7	5051	7353	0.13%	0.011%
20	8.487	927	935	944	rBV	884677	1493389	26.10%	2.190%
21	8.934	1004	1011	1024	rBV	722291	1170916	20.46%	1.717%
22	9.057	1029	1032	1036	rVV6	6151	9467	0.17%	0.014%
23	9.110	1036	1041	1049	rVB6	7719	14461	0.25%	0.021%
24	9.393	1084	1089	1096	rVB4	8119	13467	0.24%	0.020%
25	9.504	1101	1108	1114	rBV	76793	121951	2.13%	0.179%
26	9.657	1126	1134	1149	rBV	507083	896774	15.67%	1.315%
27	9.822	1156	1162	1168	rVB	3703	7924	0.14%	0.012%
28	10.204	1213	1227	1251	rBV	794179	1578614	27.58%	2.315%
29	10.581	1283	1291	1305	rBV	1644437	2843777	49.69%	4.171%
30	10.710	1305	1313	1331	rVV	874933	1596196	27.89%	2.341%
31	11.128	1377	1384	1396	rVB3	21387	51227	0.90%	0.075%
32	11.322	1411	1417	1441	rBV	135820	409125	7.15%	0.600%
33	12.081	1539	1546	1550	rVB3	6264	11314	0.20%	0.017%
34	12.169	1555	1561	1570	rBV2	16910	36972	0.65%	0.054%
35	12.410	1594	1602	1606	rBV	13254	21362	0.37%	0.031%
36	12.787	1662	1666	1673	rBV10	4613	8823	0.15%	0.013%

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37	13.169	1728	1731	1735	rBV5	3813	5747	0.10%	0.008%
38	13.222	1735	1740	1744	rVV8	4177	7379	0.13%	0.011%
39	13.263	1744	1747	1750	rVV4	7609	9717	0.17%	0.014%
40	13.328	1753	1758	1764	rVV9	10601	24315	0.42%	0.036%
41	13.398	1764	1770	1775	rVB8	7879	16345	0.29%	0.024%
42	13.622	1805	1808	1817	rBV10	6787	13545	0.24%	0.020%
43	13.834	1837	1844	1858	rBV	1569228	2405250	42.03%	3.528%
44	13.963	1864	1866	1873	rVB8	4402	8131	0.14%	0.012%
45	14.122	1882	1893	1911	rBV2	1853112	2941777	51.41%	4.315%
46	14.339	1927	1930	1935	rVB7	5846	8146	0.14%	0.012%
47	14.434	1938	1946	1965	rBV	2418650	4003701	69.96%	5.872%
48	14.639	1969	1981	1983	rBV2	1513153	2826821	49.40%	4.146%
49	14.857	2014	2018	2025	rVB10	17488	33177	0.58%	0.049%
50	14.945	2029	2033	2038	rVV7	8224	16088	0.28%	0.024%
51	15.263	2080	2087	2092	rVB5	8862	17989	0.31%	0.026%
52	15.310	2092	2095	2099	rBV6	7693	10300	0.18%	0.015%
53	15.439	2110	2117	2129	rBV	2441538	3673336	64.19%	5.387%
54	15.551	2130	2136	2154	rVV	453733	721541	12.61%	1.058%
55	15.686	2154	2159	2164	rVB7	12583	24894	0.44%	0.037%
56	15.886	2191	2193	2199	rVB7	5752	8421	0.15%	0.012%
57	16.016	2210	2215	2223	rBV5	34621	61758	1.08%	0.091%
58	16.092	2224	2228	2232	rVB7	8857	13840	0.24%	0.020%
59	16.192	2241	2245	2250	rBV6	12806	19777	0.35%	0.029%
60	16.245	2250	2254	2259	rVB7	9223	15788	0.28%	0.023%
61	16.369	2271	2275	2279	rVB7	12245	16408	0.29%	0.024%
62	16.628	2316	2319	2323	rVB6	9469	11503	0.20%	0.017%
63	16.739	2336	2338	2342	rVB5	6899	7972	0.14%	0.012%
64	16.833	2351	2354	2356	rBV4	5917	7002	0.12%	0.010%
65	17.010	2379	2384	2387	rBV5	14157	23411	0.41%	0.034%
66	17.145	2404	2407	2411	rBV6	8143	9806	0.17%	0.014%
67	17.228	2414	2421	2429	rBV	3423872	4957706	86.63%	7.271%
68	17.333	2430	2439	2451	rVV2	2542704	3975844	69.47%	5.831%
69	17.428	2451	2455	2461	rVB8	23986	38141	0.67%	0.056%
70	17.945	2538	2543	2549	rBV	62909	93309	1.63%	0.137%
71	18.051	2556	2561	2567	rBV9	22223	39370	0.69%	0.058%
72	18.104	2568	2570	2576	rBV7	9273	11903	0.21%	0.017%
73	18.169	2576	2581	2590	rBV	301561	434943	7.60%	0.638%
74	18.292	2599	2602	2608	rVB3	40604	59161	1.03%	0.087%
75	19.122	2741	2743	2744	rBV2	19381	15076	0.26%	0.022%
76	19.151	2745	2748	2753	rVB2	191169	200882	3.51%	0.295%

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Title : SVOA CALIBRATION

77	19.251	2762	2765	2769	rVB2	39497	47184	0.82%	0.069%
78	19.298	2770	2773	2774	rBV2	34216	35227	0.62%	0.052%
79	19.492	2802	2806	2812	rBV2	83935	132581	2.32%	0.194%
80	19.698	2835	2841	2847	rBV	3563721	4911344	85.82%	7.203%
81	21.351	3117	3122	3131	rBV3	249811	494334	8.64%	0.725%
82	21.680	3171	3178	3185	rBV	3297344	5450292	95.24%	7.994%
83	24.845	3708	3716	3731	rVB	1886544	4777888	83.49%	7.007%
84	25.086	3747	3757	3770	rVB	1889621	5722728	100.00%	8.393%

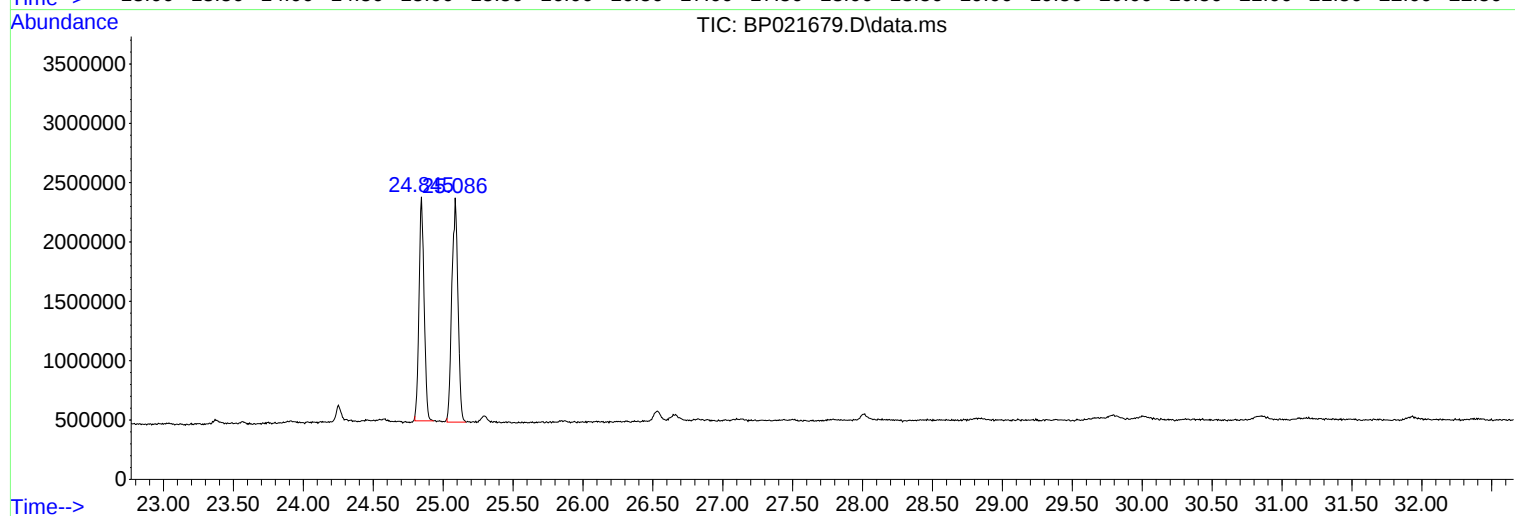
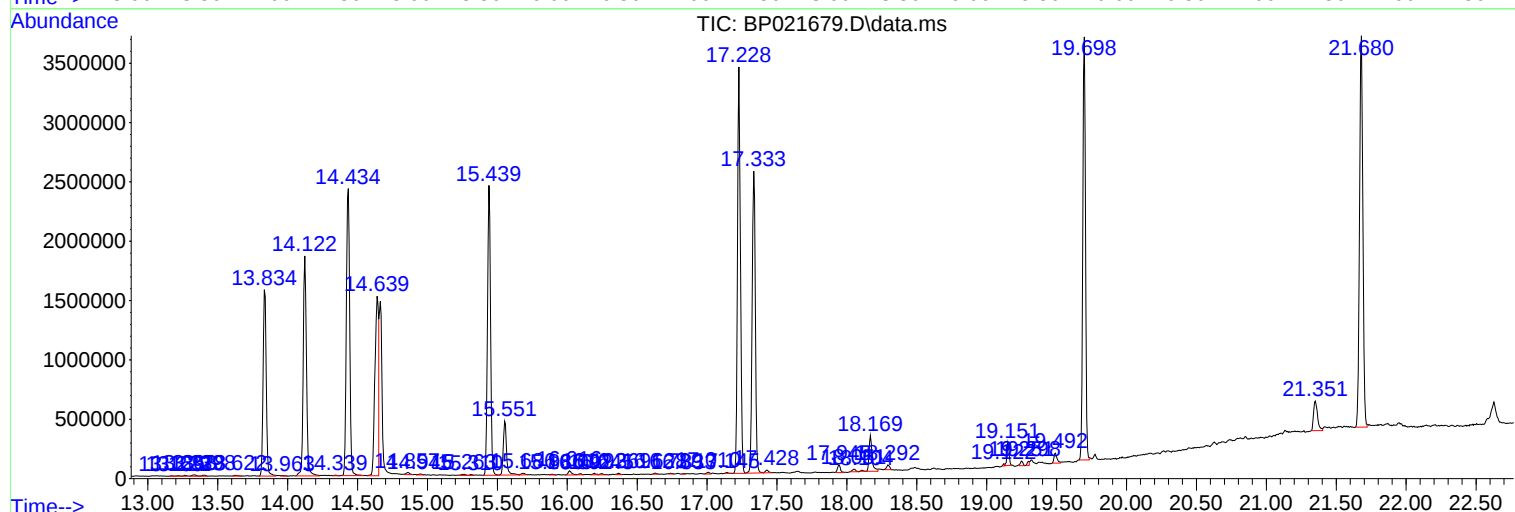
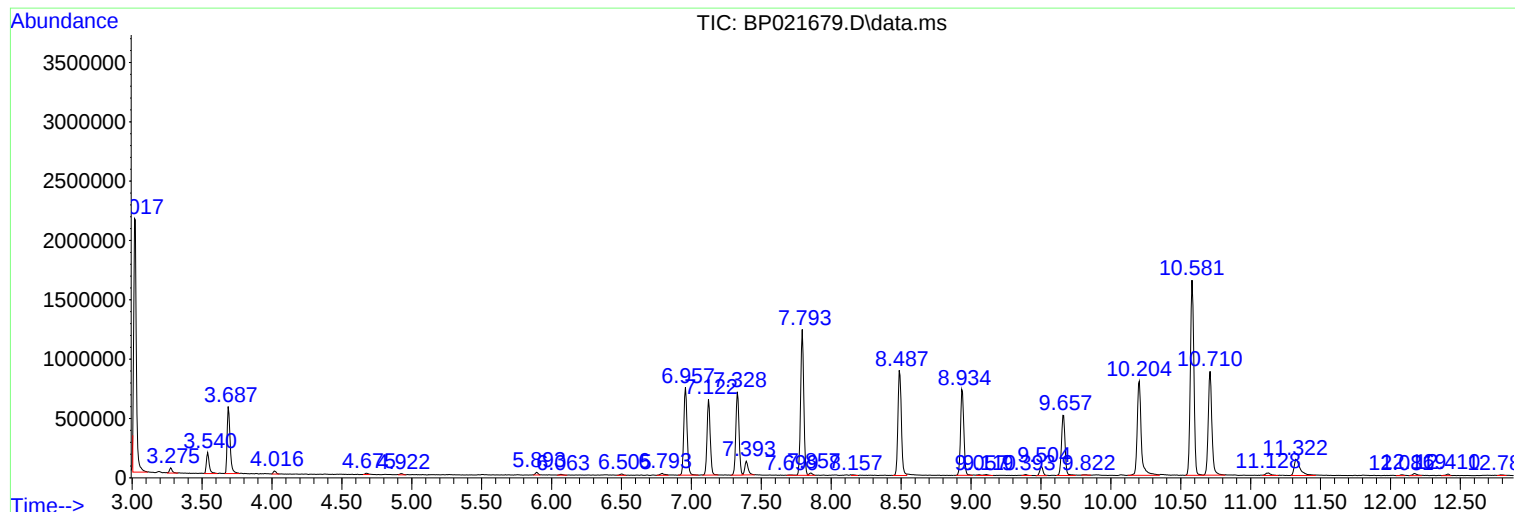
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BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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



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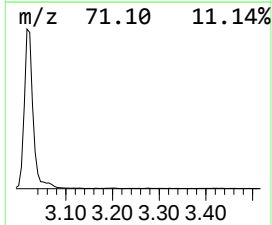
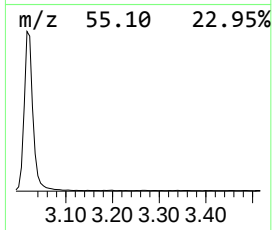
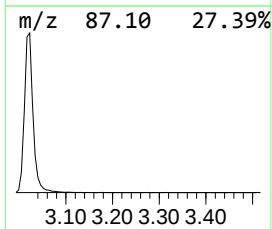
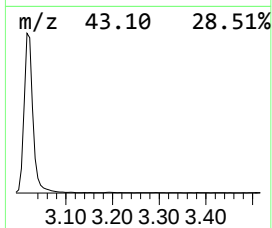
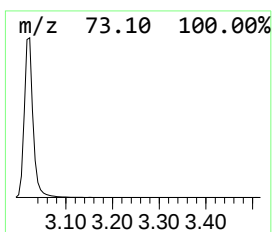
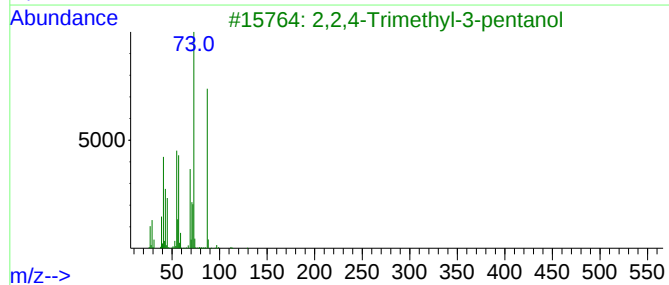
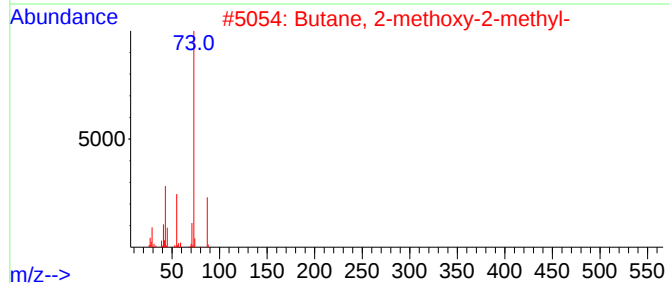
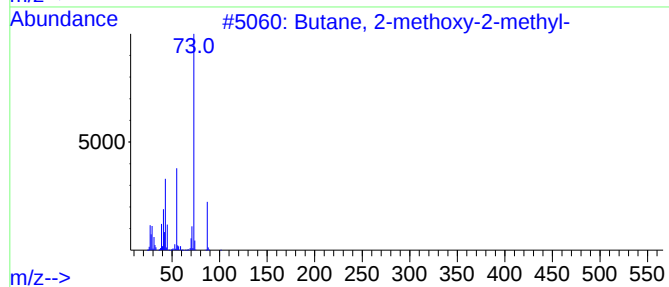
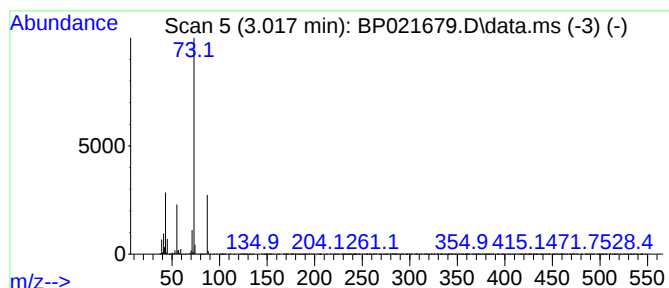
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.017	27.69 ng/ul	2690040	1,4-Dichlorobenzene-d4	7.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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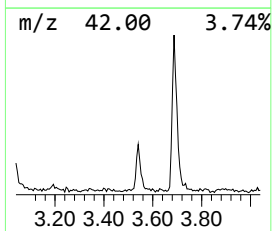
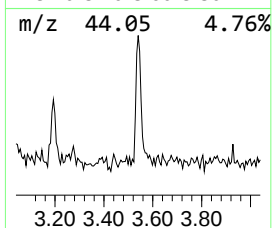
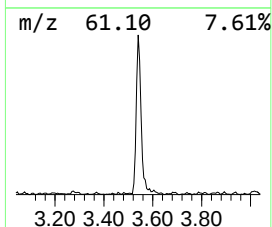
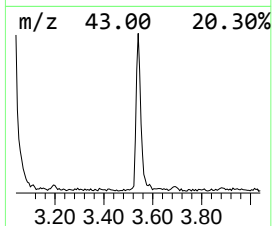
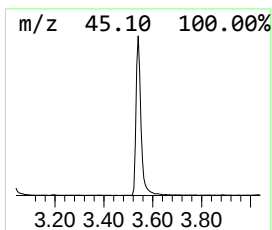
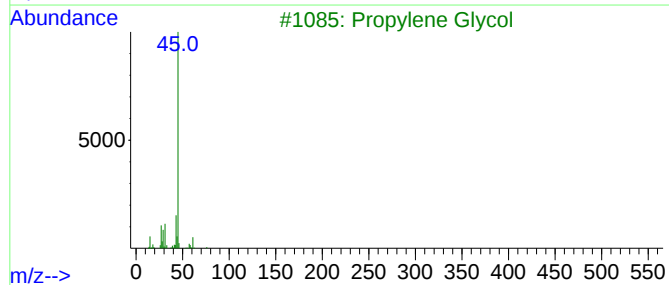
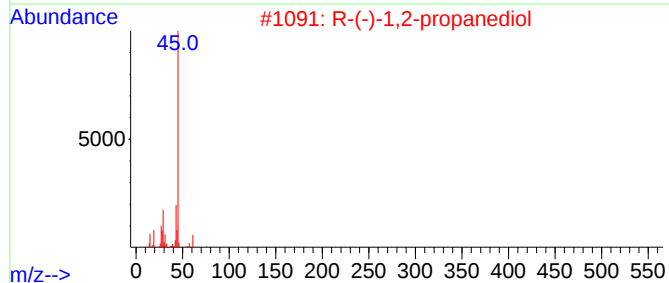
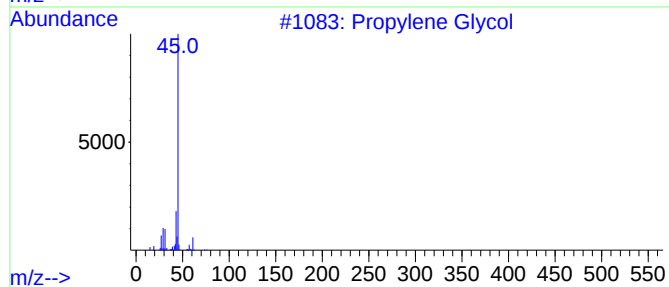
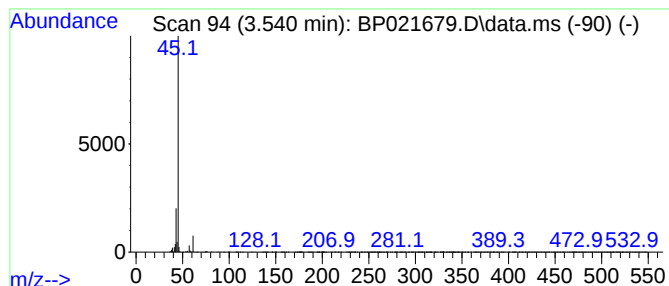
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Propylene Glycol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.540	2.52 ng/ul	244879	1,4-Dichlorobenzene-d4	7.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	90
2			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	72
3			Propylene Glycol	76	C3H8O2	000057-55-6	56
4			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
5			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	12



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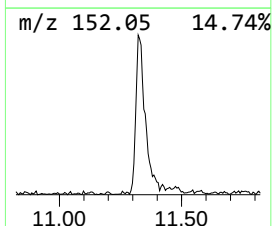
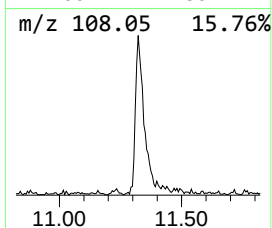
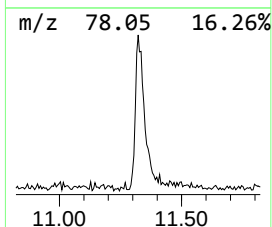
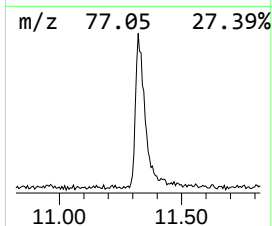
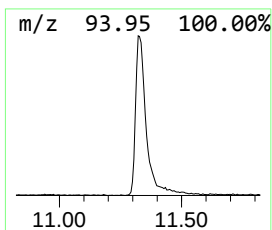
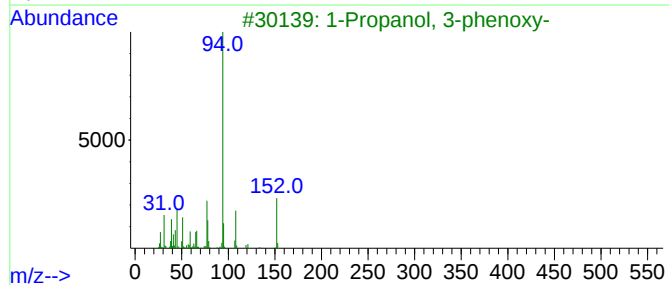
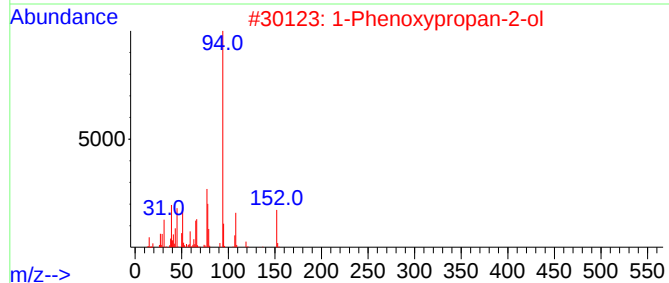
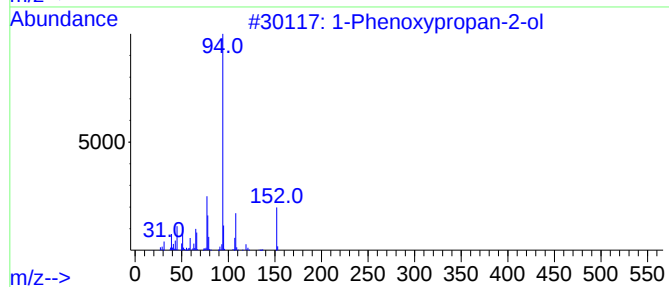
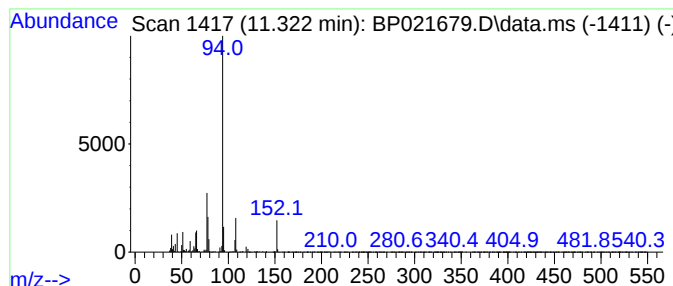
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.322	2.88 ng/ul	409125	Naphthalene-d8	10.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	97
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
3			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87



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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
Butane, 2-metho...	3.017	27.7	ng/ul	2690040	1	7.793	1942660	20.0
Propylene Glycol	3.540	2.5	ng/ul	244879	1	7.793	1942660	20.0
1-Phenoxypropan...	11.322	2.9	ng/ul	409125	2	10.581	2843780	20.0