

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083024\
 Data File : BP021759.D
 Acq On : 30 Aug 2024 19:10
 Operator : RC/JU
 Sample : P3438-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCYL1

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: BP021759.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	18	rVB	3606877	3904468	68.48%	5.490%
2	3.111	18	21	29	rVB2	29109	46156	0.81%	0.065%
3	3.269	45	48	55	rVB	51905	67173	1.18%	0.094%
4	3.540	90	94	103	rBV	186659	254079	4.46%	0.357%
5	3.687	114	119	134	rBV	660252	950896	16.68%	1.337%
6	4.011	171	174	180	rVB	19084	24135	0.42%	0.034%
7	4.675	283	287	293	rBV2	8679	14928	0.26%	0.021%
8	4.799	306	308	315	rVB8	5623	10284	0.18%	0.014%
9	4.922	324	329	339	rVB	13681	25345	0.44%	0.036%
10	5.887	489	493	499	rVB	10135	14584	0.26%	0.021%
11	6.058	517	522	530	rVB10	8506	15427	0.27%	0.022%
12	6.793	640	647	651	rBV3	10765	22381	0.39%	0.031%
13	6.957	667	675	685	rBV	788278	1310359	22.98%	1.842%
14	7.116	695	702	710	rBV	649780	1026755	18.01%	1.444%
15	7.316	729	736	743	rBV	743383	1239977	21.75%	1.743%
16	7.387	743	748	760	rVB2	61767	116372	2.04%	0.164%
17	7.781	808	815	823	rBV	1053887	1742634	30.56%	2.450%
18	7.846	823	826	830	rVB5	8369	9802	0.17%	0.014%
19	8.146	871	877	887	rVB5	7668	20633	0.36%	0.029%
20	8.487	927	935	948	rBV	1005556	1653311	29.00%	2.325%
21	8.928	1003	1010	1024	rBV	708688	1195015	20.96%	1.680%
22	9.057	1026	1032	1034	rVV6	5258	9136	0.16%	0.013%
23	9.099	1034	1039	1049	rVB6	9603	18032	0.32%	0.025%
24	9.369	1081	1085	1089	rBV7	4850	6881	0.12%	0.010%
25	9.493	1100	1106	1114	rBV	46816	74860	1.31%	0.105%
26	9.646	1123	1132	1141	rBV	556820	937320	16.44%	1.318%
27	10.193	1215	1225	1250	rBV	704511	1540475	27.02%	2.166%
28	10.398	1258	1260	1268	rVB9	5375	9177	0.16%	0.013%
29	10.569	1281	1289	1296	rBV	1308685	2369711	41.56%	3.332%
30	10.698	1303	1311	1326	rBV	885559	1545855	27.11%	2.174%
31	11.057	1367	1372	1375	rBV7	3945	6787	0.12%	0.010%
32	11.104	1375	1380	1390	rVV9	17008	41421	0.73%	0.058%
33	11.310	1406	1415	1432	rBV	138260	410109	7.19%	0.577%
34	12.163	1551	1560	1566	rBV4	13671	29828	0.52%	0.042%
35	12.234	1568	1572	1577	rVB7	3907	6694	0.12%	0.009%
36	12.398	1596	1600	1605	rVV2	6314	9046	0.16%	0.013%

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Integrator: RTE

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

37	12.457	1605	1610	1615	rVB4	10199	19029	0.33%	0.027%
38	12.775	1661	1664	1672	rBV10	4217	9133	0.16%	0.013%
39	12.904	1681	1686	1690	rBV8	3629	6233	0.11%	0.009%
40	13.034	1703	1708	1710	rBV5	4928	9171	0.16%	0.013%
41	13.257	1741	1746	1751	rVB8	6237	11337	0.20%	0.016%
42	13.322	1751	1757	1763	rBV6	10201	19734	0.35%	0.028%
43	13.387	1763	1768	1771	rVV7	7331	13334	0.23%	0.019%
44	13.445	1773	1778	1781	rVV6	4657	10315	0.18%	0.015%
45	13.604	1795	1805	1806	rBV4	16983	33575	0.59%	0.047%
46	13.745	1823	1829	1835	rVB3	16585	28808	0.51%	0.041%
47	13.828	1835	1843	1857	rBV	1661238	2404651	42.17%	3.381%
48	13.981	1864	1869	1873	rBV6	9838	16346	0.29%	0.023%
49	14.116	1881	1892	1909	rBV	1784488	2823292	49.52%	3.970%
50	14.334	1926	1929	1934	rVB6	6039	7666	0.13%	0.011%
51	14.422	1934	1944	1951	rBV	2014792	3222283	56.51%	4.531%
52	14.492	1951	1956	1965	rVB	246831	413414	7.25%	0.581%
53	14.645	1974	1982	1983	rBV3	2297189	3656245	64.12%	5.141%
54	14.834	2008	2014	2024	rVB	168587	298689	5.24%	0.420%
55	14.916	2024	2028	2030	rVV4	6255	8935	0.16%	0.013%
56	14.945	2030	2033	2036	rVB4	11555	13370	0.23%	0.019%
57	15.039	2044	2049	2050	rBV5	6167	9373	0.16%	0.013%
58	15.057	2050	2052	2054	rVV3	8061	8402	0.15%	0.012%
59	15.110	2058	2061	2066	rVB7	9541	13898	0.24%	0.020%
60	15.245	2077	2084	2091	rBV7	10427	27870	0.49%	0.039%
61	15.310	2091	2095	2099	rVB6	6279	8925	0.16%	0.013%
62	15.434	2109	2116	2122	rBV	2477459	3644633	63.92%	5.125%
63	15.492	2122	2126	2131	rVB	171925	264896	4.65%	0.372%
64	15.551	2131	2136	2144	rBV	652321	818184	14.35%	1.150%
65	15.657	2150	2154	2161	rVB3	28759	45879	0.80%	0.065%
66	15.745	2165	2169	2171	rBV4	9638	11407	0.20%	0.016%
67	15.792	2173	2177	2187	rVB	37793	66164	1.16%	0.093%
68	15.886	2187	2193	2196	rBV8	6097	11605	0.20%	0.016%
69	15.939	2198	2202	2210	rVB	38440	71062	1.25%	0.100%
70	16.004	2210	2213	2225	rVB9	23433	49227	0.86%	0.069%
71	16.181	2233	2243	2248	rBV9	18695	49835	0.87%	0.070%
72	16.304	2260	2264	2268	rVB7	5133	8052	0.14%	0.011%
73	16.492	2290	2296	2297	rBV6	8255	15140	0.27%	0.021%
74	16.516	2297	2300	2305	rVV5	17607	27982	0.49%	0.039%
75	16.563	2306	2308	2314	rVB5	7966	10364	0.18%	0.015%
76	16.622	2315	2318	2320	rBV4	6914	7250	0.13%	0.010%

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77	16.951	2370	2374	2377	rBV6	12355	18993	0.33%	0.027%
78	17.039	2385	2389	2397	rVB	49389	90915	1.59%	0.128%
79	17.228	2414	2421	2425	rBV	2643127	4092184	71.77%	5.754%
80	17.269	2425	2428	2432	rVV	505180	674873	11.84%	0.949%
81	17.328	2432	2438	2450	rVV2	2656934	4213855	73.90%	5.925%
82	17.422	2451	2454	2459	rVB4	35013	47853	0.84%	0.067%
83	17.645	2489	2492	2497	rVB4	14769	22062	0.39%	0.031%
84	17.763	2510	2512	2515	rBV4	8447	8453	0.15%	0.012%
85	17.939	2536	2542	2547	rBV	53724	80680	1.41%	0.113%
86	18.127	2570	2574	2576	rBV	37419	50535	0.89%	0.071%
87	18.163	2576	2580	2589	rVB2	186631	329543	5.78%	0.463%
88	18.269	2595	2598	2605	rBV6	21390	38833	0.68%	0.055%
89	18.339	2605	2610	2619	rVB2	84865	161302	2.83%	0.227%
90	18.651	2659	2663	2667	rBV3	37452	58757	1.03%	0.083%
91	19.151	2745	2748	2756	rVB7	53274	76452	1.34%	0.107%
92	19.251	2762	2765	2769	rVB3	39130	49269	0.86%	0.069%
93	19.298	2770	2773	2774	rBV3	21673	25372	0.44%	0.036%
94	19.351	2778	2782	2787	rVB	531836	672969	11.80%	0.946%
95	19.486	2803	2805	2811	rVB5	38091	45576	0.80%	0.064%
96	19.698	2835	2841	2850	rBV2	3932560	5701836	100.00%	8.017%
97	21.357	3118	3123	3131	rVB	248599	439408	7.71%	0.618%
98	21.680	3172	3178	3185	rBV2	2872603	4859610	85.23%	6.833%
99	24.857	3709	3718	3733	rVB	1728860	5329472	93.47%	7.493%
100	25.092	3748	3758	3771	rVB2	1678958	5156850	90.44%	7.251%

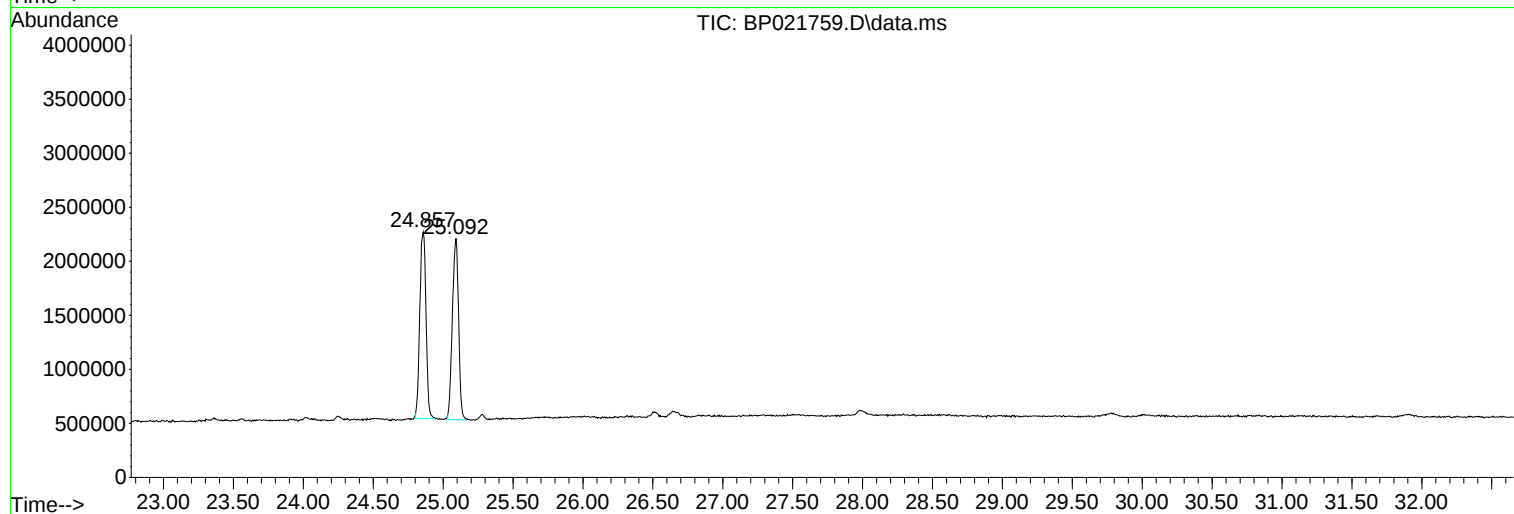
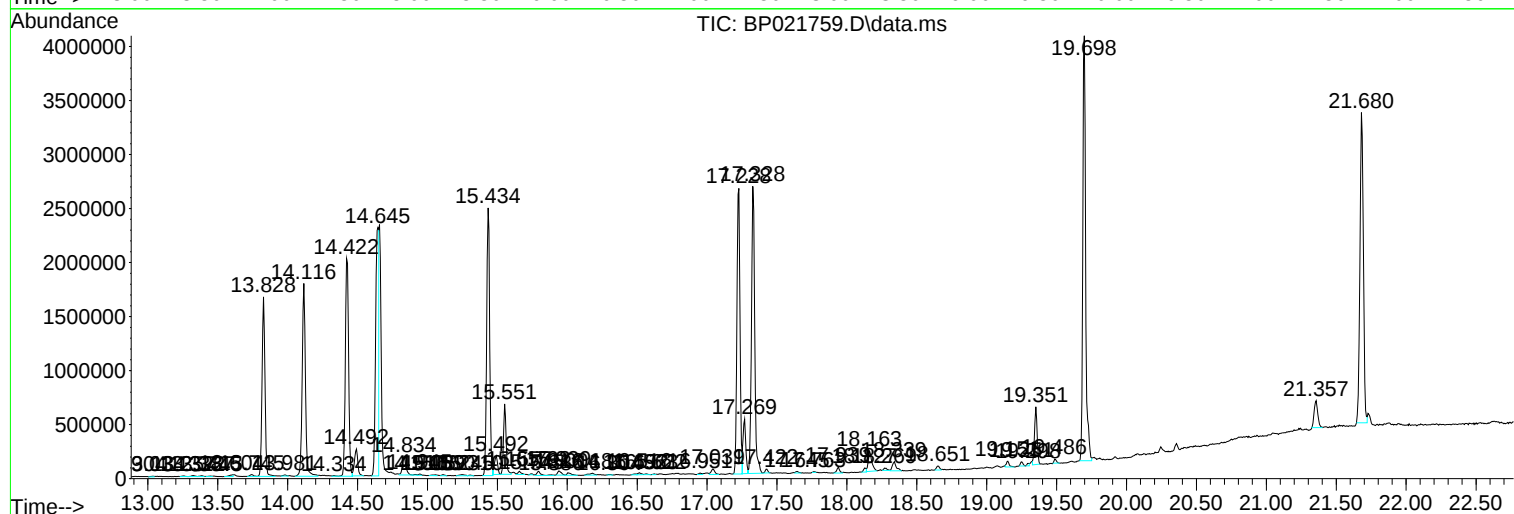
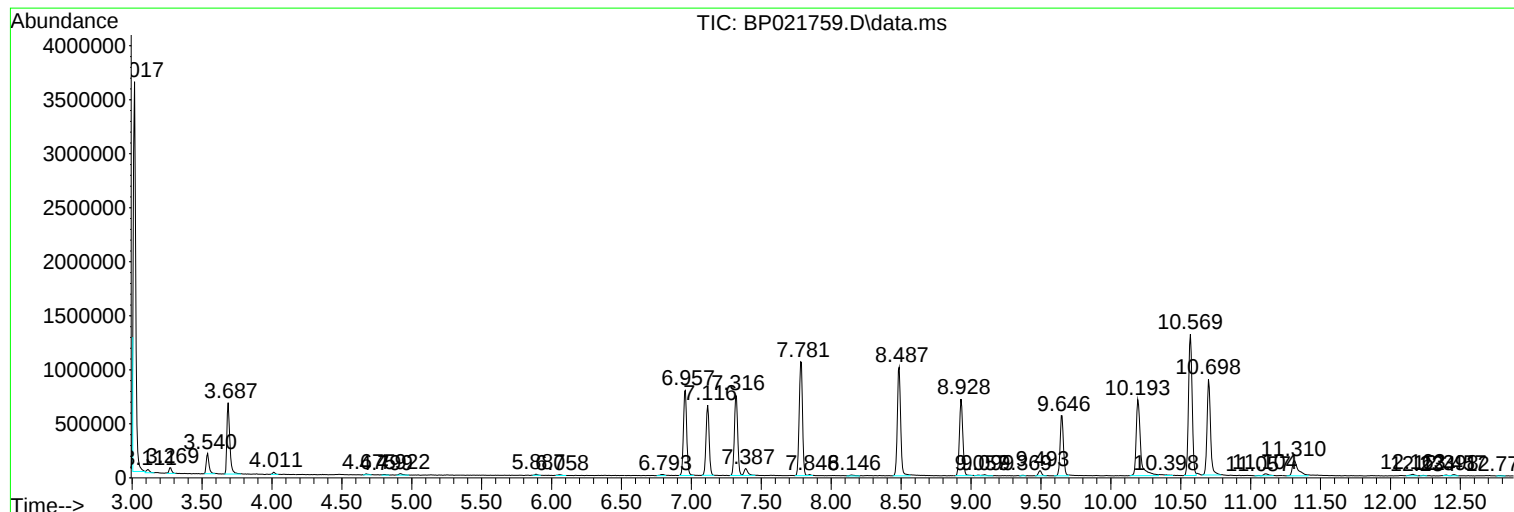
Sum of corrected areas: 71121406

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083024\
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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYL1

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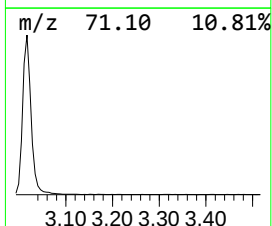
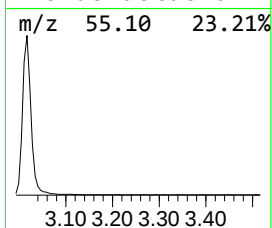
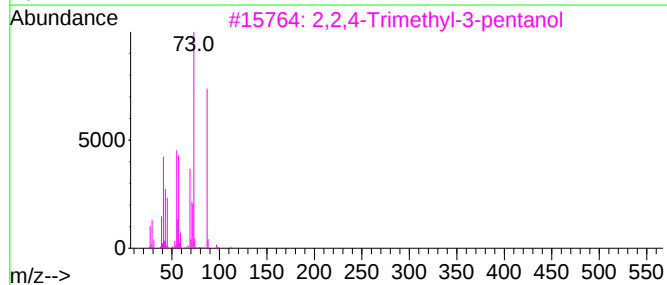
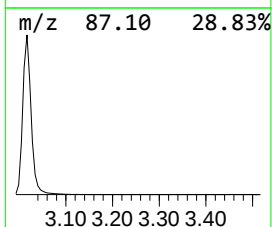
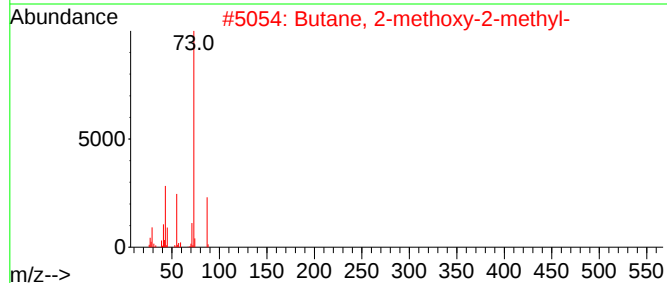
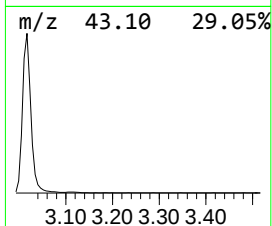
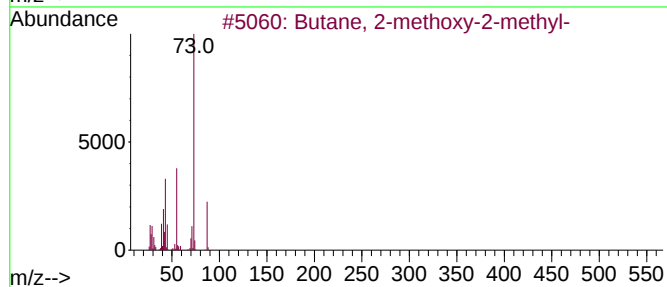
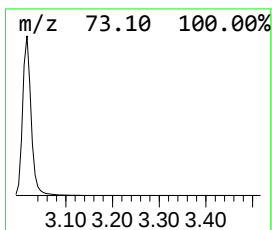
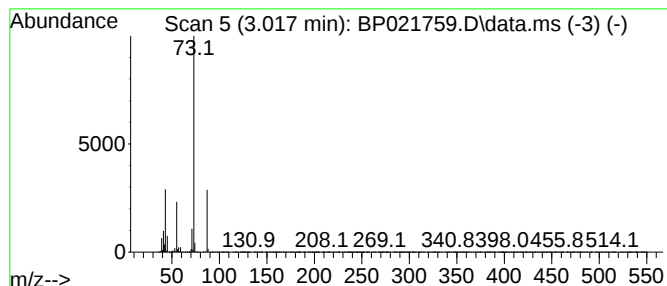
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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	44.81 ng/ul	3904470	1,4-Dichlorobenzene-d4	7.787

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78		
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39		
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23		
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9		



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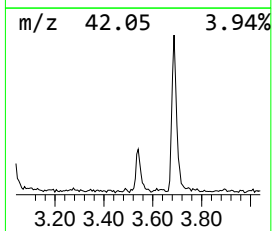
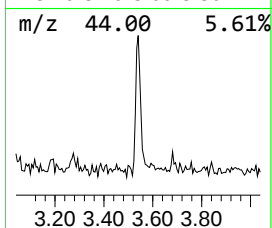
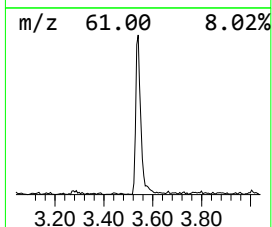
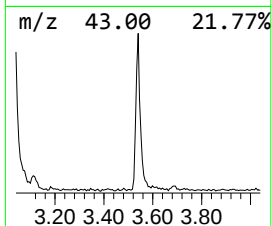
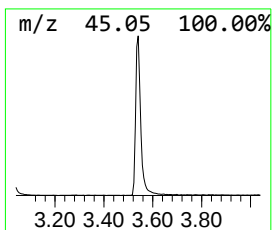
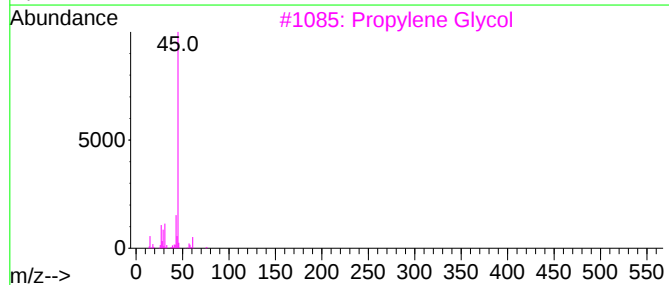
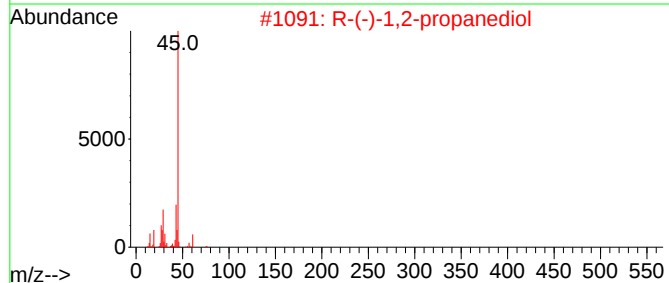
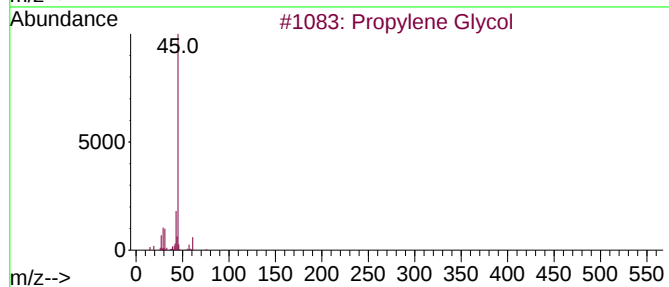
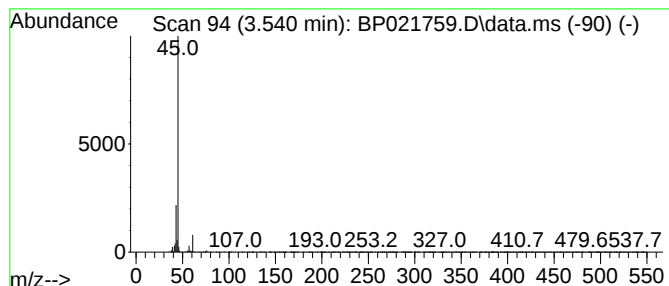
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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	2.92 ng/ul	254079	1,4-Dichlorobenzene-d4	7.787

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	78
2			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	78
3			Propylene Glycol	76	C3H8O2	000057-55-6	56
4			Propylene Glycol	76	C3H8O2	000057-55-6	56
5			Isopropyl Alcohol	60	C3H8O	000067-63-0	56



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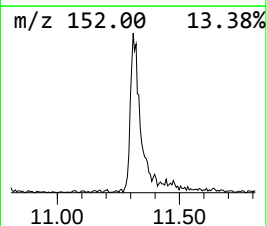
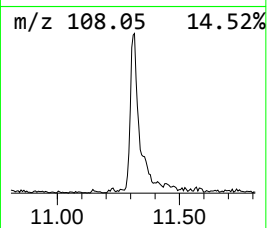
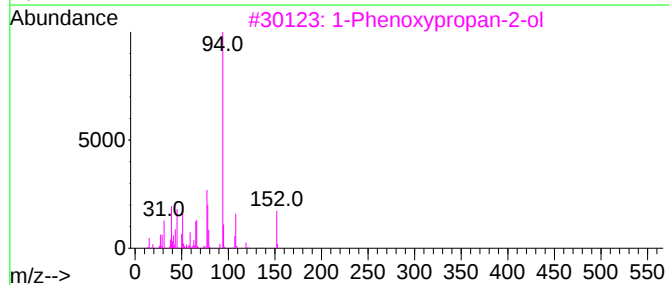
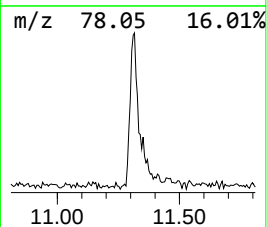
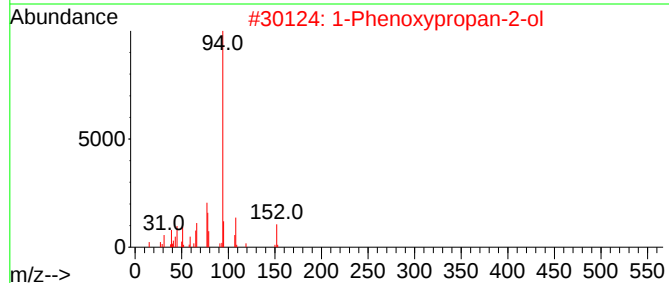
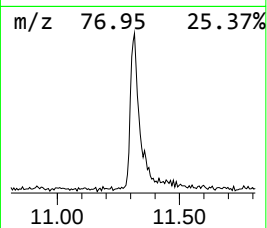
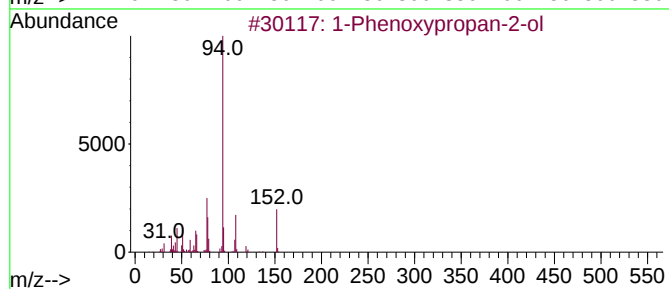
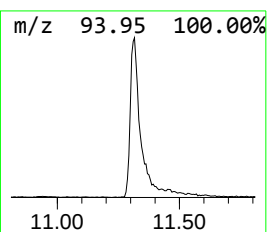
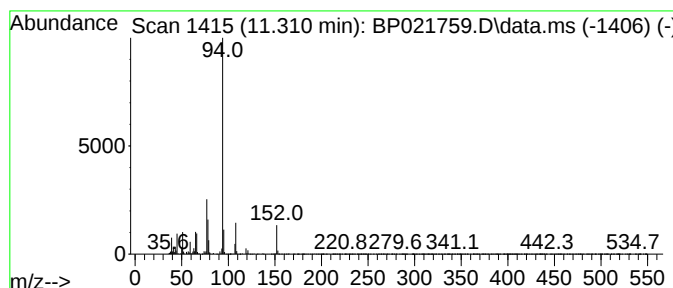
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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.
11.310	3.46 ng/ul	410109	Naphthalene-d8	10.569

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	96
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	91
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083024\
Data File : BP021759.D
Acq On : 30 Aug 2024 19:10
Operator : RC/JU
Sample : P3438-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYL1

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.017	44.8	ng/ul	3904470	1	7.787	1742630	20.0
Propylene Glycol	3.540	2.9	ng/ul	254079	1	7.787	1742630	20.0
1-Phenoxypropan...	11.310	3.5	ng/ul	410109	2	10.569	2369710	20.0