

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
 Data File : BP021678.D
 Acq On : 27 Aug 2024 19:43
 Operator : RC/JU
 Sample : P3675-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCYC4

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: BP021678.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.022	3	6	26	rVB	4150297	5149800	84.72%	5.989%
2	3.275	45	49	56	rVB	72948	98723	1.62%	0.115%
3	3.540	90	94	102	rBV	216815	292438	4.81%	0.340%
4	3.687	115	119	138	rVB	920225	1305592	21.48%	1.518%
5	4.022	172	176	181	rVB	26659	37609	0.62%	0.044%
6	4.681	284	288	290	rBV4	8557	11936	0.20%	0.014%
7	4.805	304	309	315	rBV10	5484	10957	0.18%	0.013%
8	4.928	325	330	336	rVB	13342	22076	0.36%	0.026%
9	5.893	489	494	501	rVB	36406	55500	0.91%	0.065%
10	6.063	517	523	529	rVB10	9571	16961	0.28%	0.020%
11	6.499	592	597	604	rVB2	22545	36957	0.61%	0.043%
12	6.793	641	647	654	rBV2	16256	36081	0.59%	0.042%
13	6.958	666	675	689	rBV	1112736	1778956	29.27%	2.069%
14	7.122	697	703	712	rVB	936373	1446098	23.79%	1.682%
15	7.246	720	724	730	rVB6	7208	11307	0.19%	0.013%
16	7.328	730	738	744	rBV	1024583	1686388	27.74%	1.961%
17	7.393	744	749	759	rVB	126818	213989	3.52%	0.249%
18	7.793	809	817	824	rBV	1298000	2073216	34.11%	2.411%
19	7.852	824	827	835	rVV3	29296	43897	0.72%	0.051%
20	8.105	865	870	871	rBV4	4572	7611	0.13%	0.009%
21	8.152	874	878	882	rVB6	4572	9003	0.15%	0.010%
22	8.487	928	935	944	rBV	1337760	2199551	36.18%	2.558%
23	8.940	1004	1012	1023	rBV	986335	1624920	26.73%	1.890%
24	9.052	1028	1031	1037	rVV7	6153	12270	0.20%	0.014%
25	9.110	1037	1041	1045	rVB7	8064	12587	0.21%	0.015%
26	9.387	1079	1088	1092	rBV4	13944	23144	0.38%	0.027%
27	9.499	1101	1107	1114	rBV	99970	149361	2.46%	0.174%
28	9.657	1127	1134	1142	rBV	768035	1222460	20.11%	1.422%
29	9.893	1168	1174	1176	rBV7	8855	14774	0.24%	0.017%
30	10.199	1217	1226	1244	rBV	1098583	2078928	34.20%	2.418%
31	10.581	1282	1291	1298	rBV	1587747	2813713	46.29%	3.272%
32	10.704	1305	1312	1331	rBV	1130748	2079228	34.20%	2.418%
33	11.116	1373	1382	1391	rBV5	26698	59963	0.99%	0.070%
34	11.322	1409	1417	1430	rBV	174948	443132	7.29%	0.515%
35	12.010	1528	1534	1538	rBV8	4932	9392	0.15%	0.011%
36	12.169	1553	1561	1570	rBV2	23031	49561	0.82%	0.058%

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37	12.404	1597	1601	1607	rVB2	7234	13924	0.23%	0.016%
38	12.798	1662	1668	1677	rVV3	168690	274281	4.51%	0.319%
39	13.040	1703	1709	1711	rBV5	4546	7725	0.13%	0.009%
40	13.122	1717	1723	1727	rVB9	5132	10646	0.18%	0.012%
41	13.175	1727	1732	1734	rBV4	42725	69905	1.15%	0.081%
42	13.198	1734	1736	1743	rVB3	52507	78699	1.29%	0.092%
43	13.263	1743	1747	1750	rBV5	11788	17381	0.29%	0.020%
44	13.316	1750	1756	1763	rVB	181190	280835	4.62%	0.327%
45	13.392	1765	1769	1773	rVB6	6914	12387	0.20%	0.014%
46	13.481	1780	1784	1787	rBV4	4487	8676	0.14%	0.010%
47	13.716	1819	1824	1825	rBV5	5730	8512	0.14%	0.010%
48	13.763	1825	1832	1838	rVV	1329937	1833310	30.16%	2.132%
49	13.834	1838	1844	1859	rVV	2007296	3282853	54.01%	3.818%
50	13.957	1863	1865	1875	rVB	5653	11379	0.19%	0.013%
51	14.122	1877	1893	1914	rBV	2360034	3761570	61.88%	4.374%
52	14.287	1914	1921	1926	rBV	85522	130639	2.15%	0.152%
53	14.345	1926	1931	1933	rBV6	9333	15138	0.25%	0.018%
54	14.434	1935	1946	1955	rVV	2379207	3821771	62.87%	4.444%
55	14.516	1955	1960	1970	rVB	110906	193659	3.19%	0.225%
56	14.628	1970	1979	1988	rBV	791549	1767418	29.08%	2.055%
57	14.698	1988	1991	1997	rVV	242551	370535	6.10%	0.431%
58	14.769	1997	2003	2011	rVB2	97102	204371	3.36%	0.238%
59	14.857	2014	2018	2025	rVB7	25383	53605	0.88%	0.062%
60	14.951	2030	2034	2035	rVV4	5441	7833	0.13%	0.009%
61	14.975	2035	2038	2044	rVB4	17027	27726	0.46%	0.032%
62	15.045	2046	2050	2053	rBV6	4938	9501	0.16%	0.011%
63	15.175	2068	2072	2077	rVB2	19033	27141	0.45%	0.032%
64	15.251	2081	2085	2089	rBV5	10670	17475	0.29%	0.020%
65	15.316	2091	2096	2098	rBV4	12799	22054	0.36%	0.026%
66	15.339	2098	2100	2102	rVB3	9450	7953	0.13%	0.009%
67	15.439	2110	2117	2129	rBV	3207792	4669728	76.82%	5.430%
68	15.551	2130	2136	2144	rBV	640343	983715	16.18%	1.144%
69	15.681	2155	2158	2166	rVB9	15255	30917	0.51%	0.036%
70	15.792	2174	2177	2188	rVB9	22335	44937	0.74%	0.052%
71	15.892	2188	2194	2195	rBV5	9998	16182	0.27%	0.019%
72	15.934	2195	2201	2207	rVB6	35171	77766	1.28%	0.090%
73	16.016	2210	2215	2216	rBV3	34535	46649	0.77%	0.054%
74	16.110	2227	2231	2236	rVB	15735	20892	0.34%	0.024%
75	16.186	2239	2244	2250	rVB10	19525	37891	0.62%	0.044%
76	16.322	2262	2267	2269	rBV5	11939	18876	0.31%	0.022%

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 Title : SVOA CALIBRATION

77	16.363	2272	2274	2283	rVB10	14273	24891	0.41%	0.029%
78	16.733	2334	2337	2343	rVB8	10102	18784	0.31%	0.022%
79	16.792	2343	2347	2349	rBV5	10782	17294	0.28%	0.020%
80	16.951	2371	2374	2376	rBV4	6104	8681	0.14%	0.010%
81	17.004	2379	2383	2387	rBV4	19082	29742	0.49%	0.035%
82	17.175	2407	2412	2413	rBV5	10564	15177	0.25%	0.018%
83	17.228	2414	2421	2430	rVV	3175837	4504377	74.10%	5.238%
84	17.328	2431	2438	2449	rVV2	2928800	4739919	77.97%	5.512%
85	17.422	2451	2454	2459	rVB6	27221	45652	0.75%	0.053%
86	17.945	2538	2543	2547	rVB	62809	93002	1.53%	0.108%
87	18.169	2576	2581	2587	rBV	278533	422105	6.94%	0.491%
88	18.292	2598	2602	2608	rVB4	50747	69869	1.15%	0.081%
89	18.586	2649	2652	2655	rBV3	22171	28398	0.47%	0.033%
90	19.098	2727	2739	2745	rBV5	278049	984585	16.20%	1.145%
91	19.151	2745	2748	2752	rVB4	140175	203209	3.34%	0.236%
92	19.233	2759	2762	2763	rBV2	58820	64677	1.06%	0.075%
93	19.492	2803	2806	2814	rBV4	92001	146260	2.41%	0.170%
94	19.698	2835	2841	2850	rVB	4009326	5845989	96.17%	6.798%
95	20.080	2897	2906	2908	rBV2	442559	931406	15.32%	1.083%
96	20.186	2918	2924	2935	rVB5	394606	1174296	19.32%	1.366%
97	21.257	3101	3106	3107	rBV4	232431	365247	6.01%	0.425%
98	21.674	3171	3177	3187	rBV2	3079177	5198381	85.52%	6.045%
99	24.851	3707	3717	3729	rVB	2290054	6078774	100.00%	7.069%
100	25.086	3747	3757	3772	rVB	1871980	5545746	91.23%	6.449%

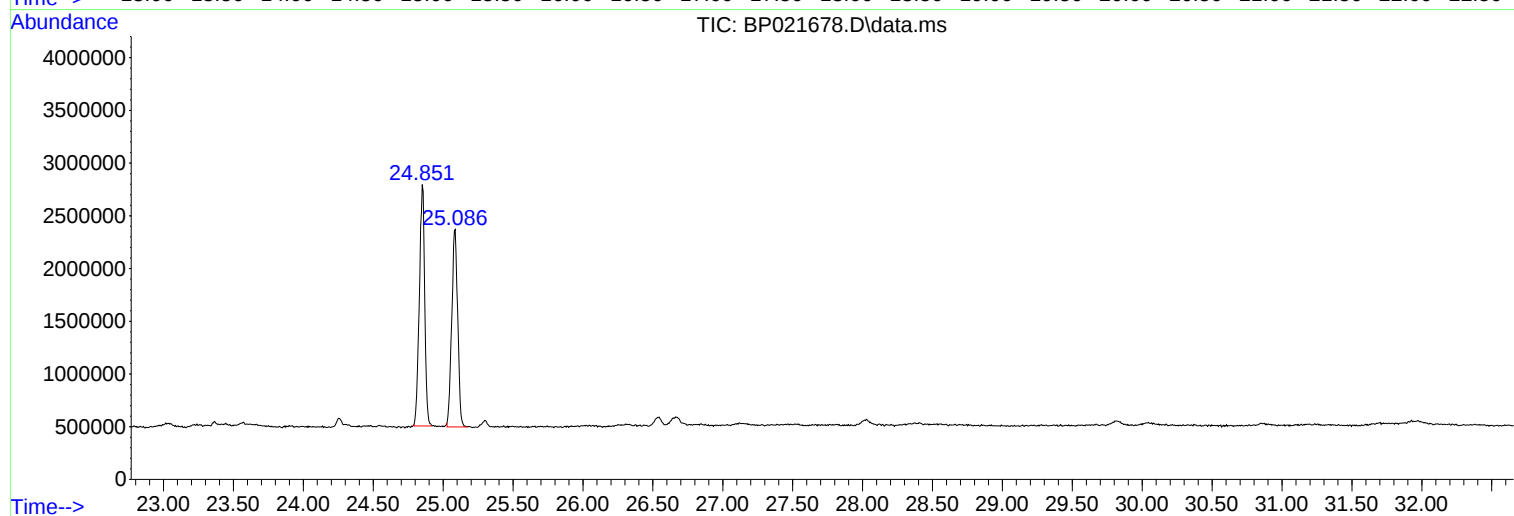
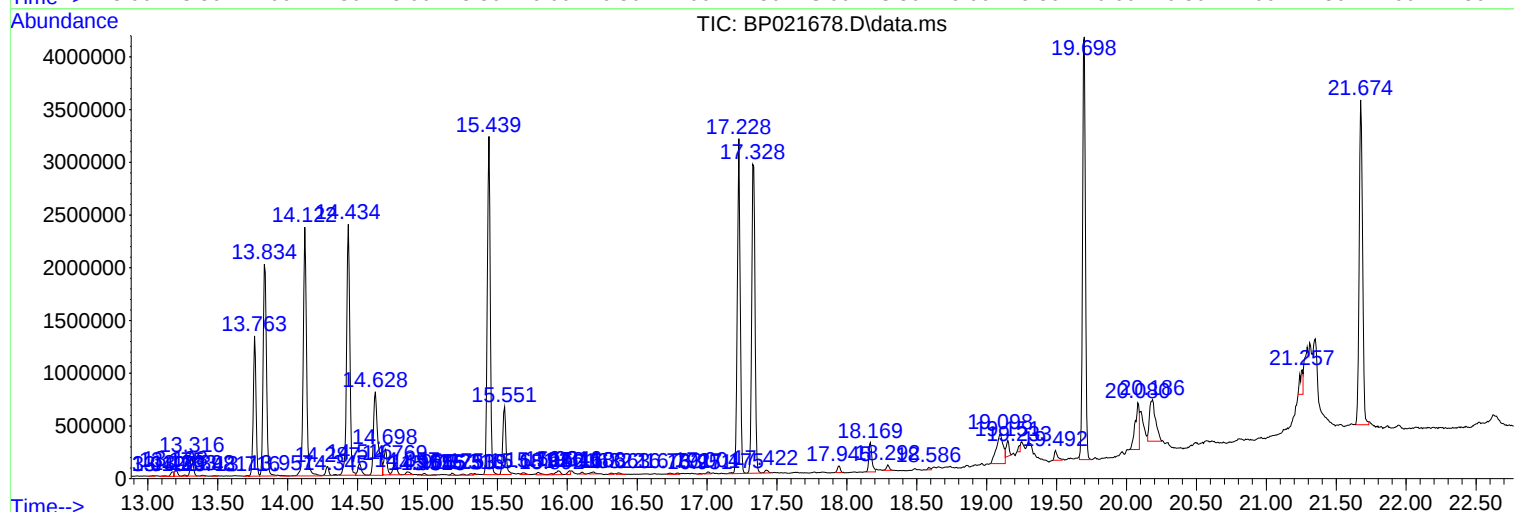
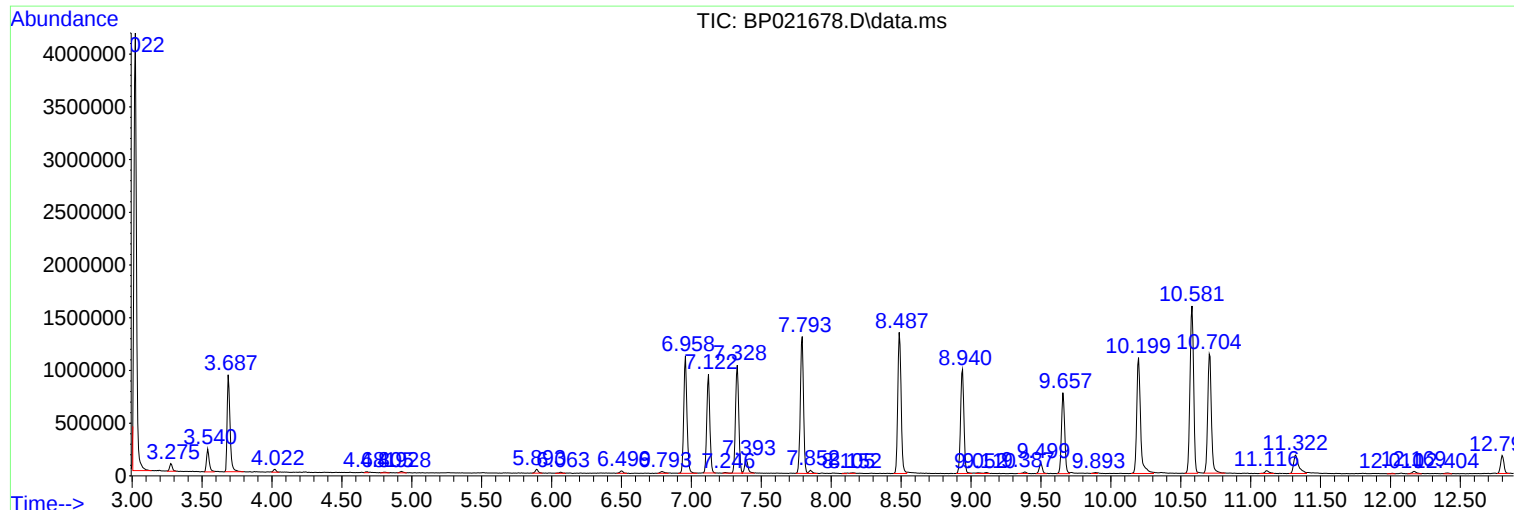
Sum of corrected areas: 85990995

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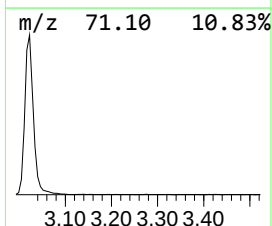
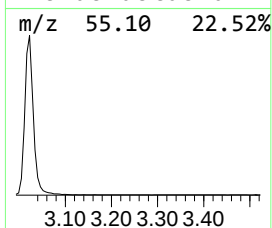
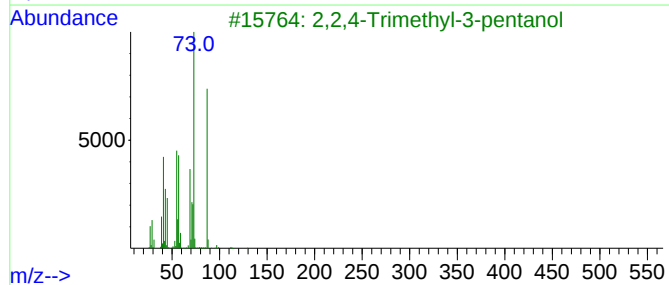
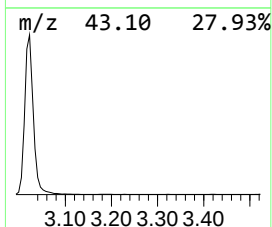
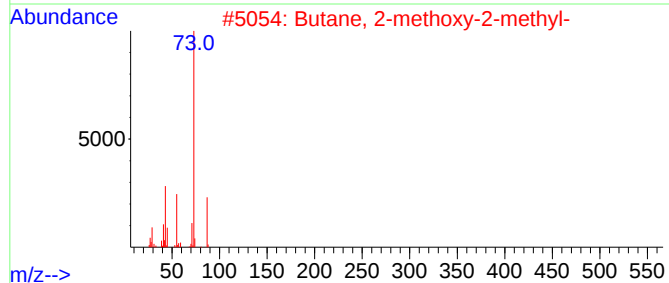
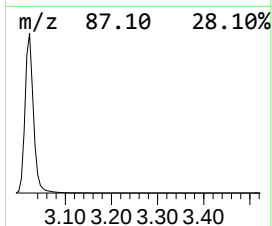
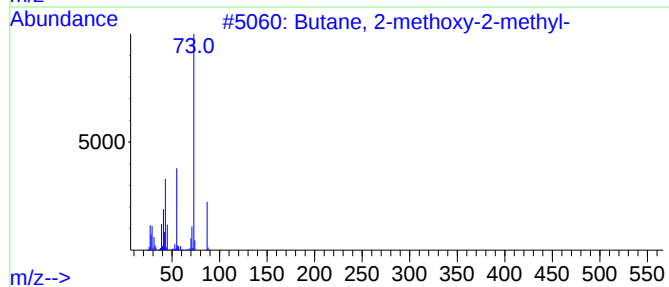
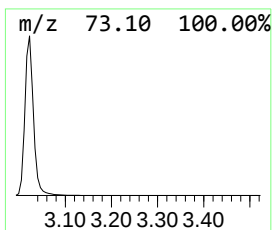
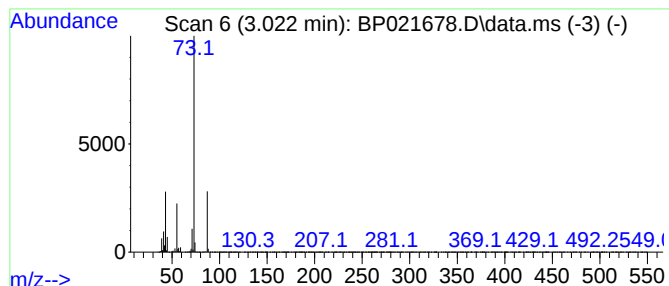
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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.
3.022	49.68 ng/ul	5149800	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	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9		



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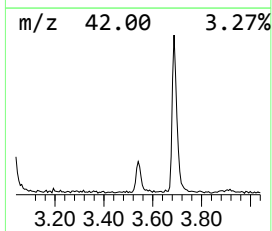
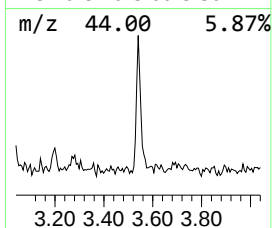
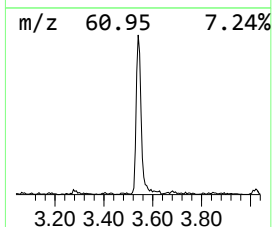
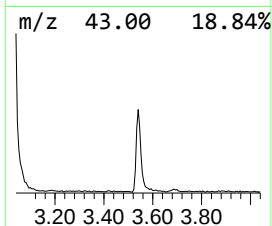
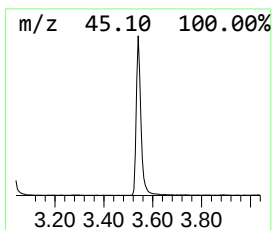
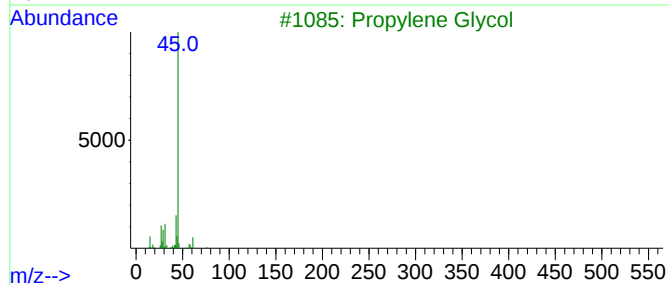
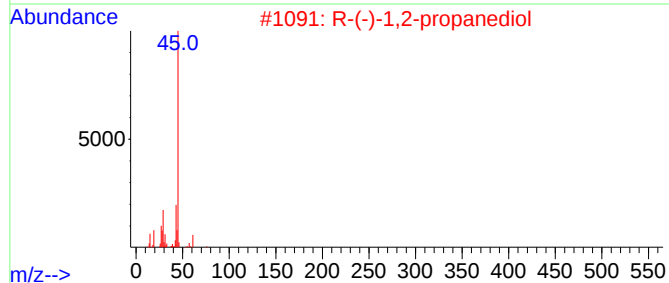
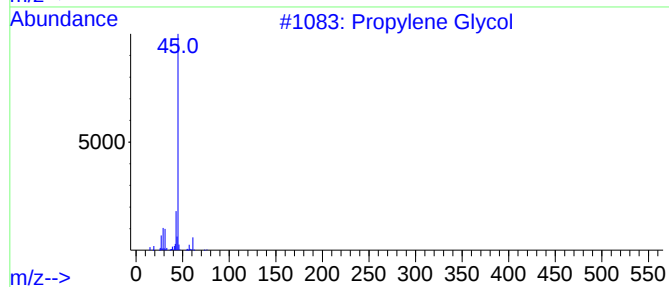
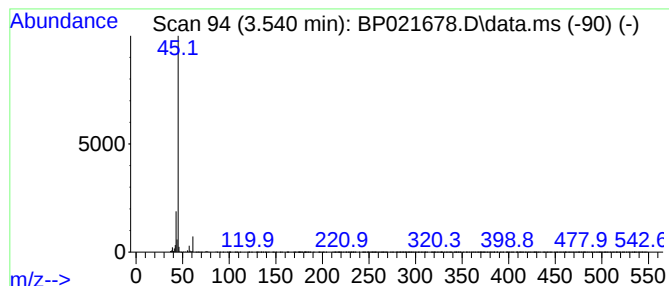
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Propylene Glycol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.540	2.82 ng/ul	292438	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	64
4			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
5			2-Butanol	74	C4H10O	000078-92-2	40



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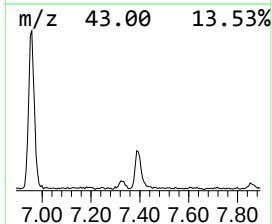
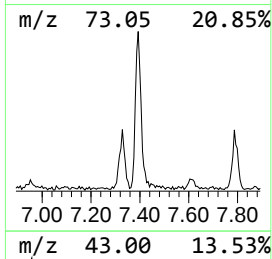
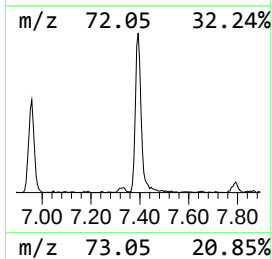
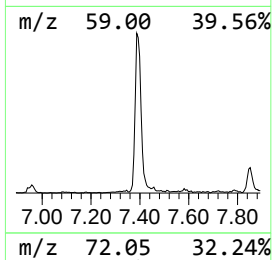
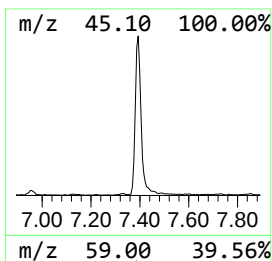
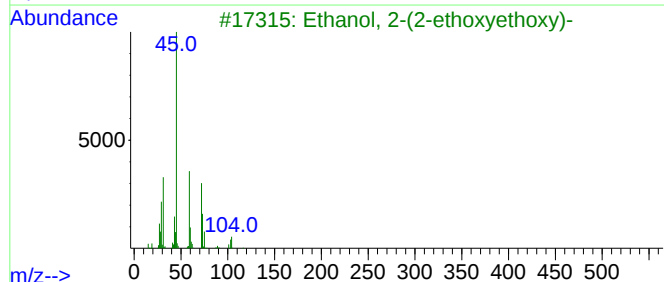
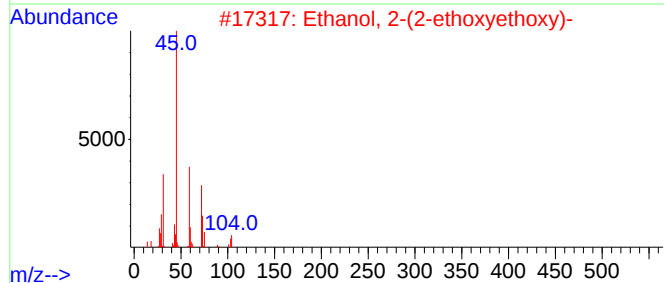
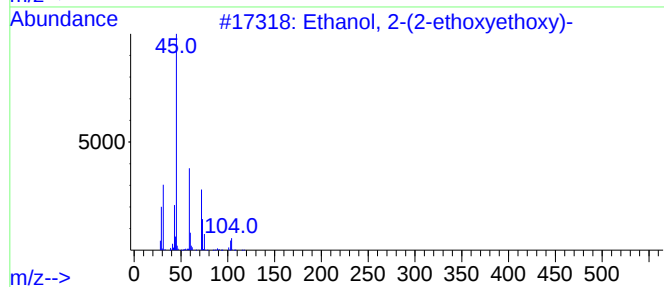
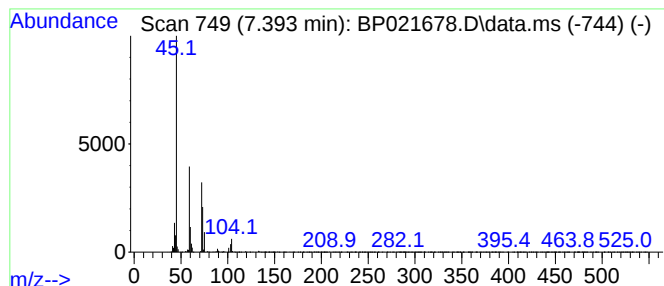
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.393	2.06 ng/ul	213989	1,4-Dichlorobenzene-d4	7.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	86
2			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	86
3			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	86
4			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	78
5			Diethyl carbitol	162	C8H18O3	000112-36-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021678.D
Acq On : 27 Aug 2024 19:43
Operator : RC/JU
Sample : P3675-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYC4

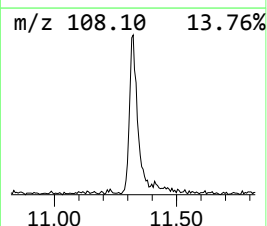
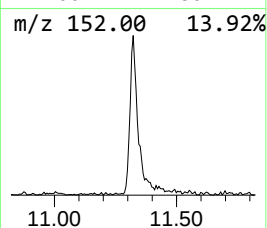
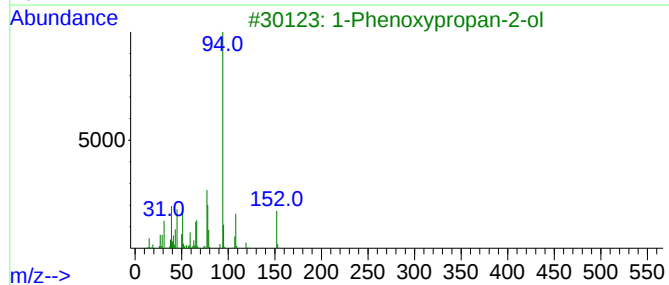
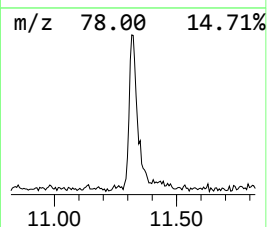
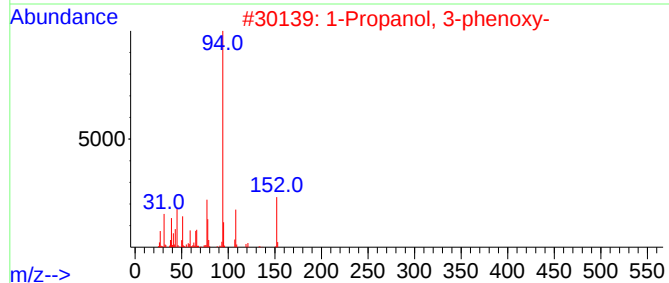
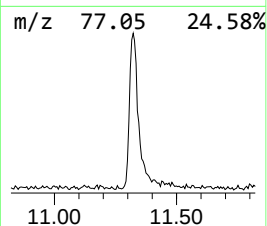
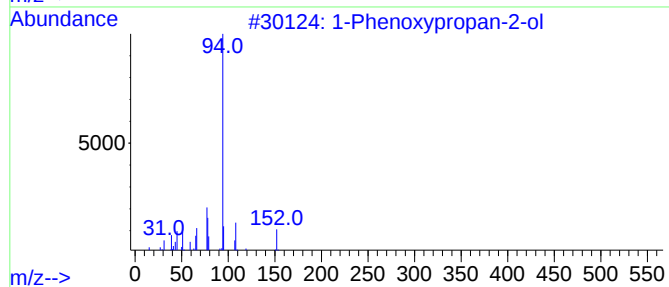
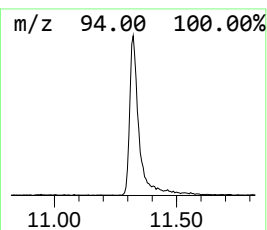
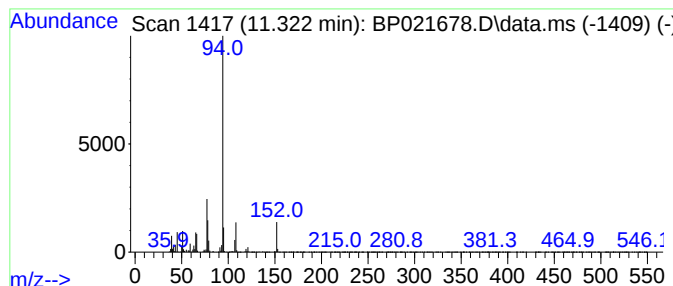
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.322	3.15 ng/ul	443132	Naphthalene-d8	10.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
2			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	83
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	70
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	70
5			Carbonic acid, neopentyl phenyl ...	208	C12H16O3	013183-19-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021678.D
Acq On : 27 Aug 2024 19:43
Operator : RC/JU
Sample : P3675-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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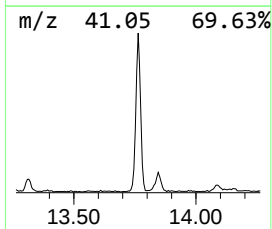
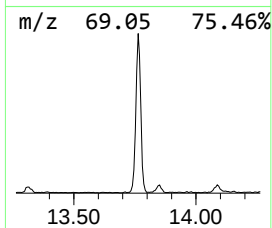
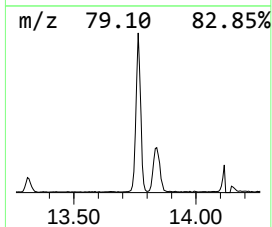
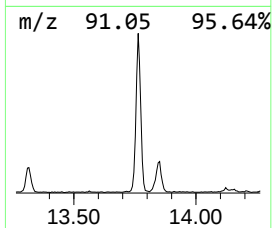
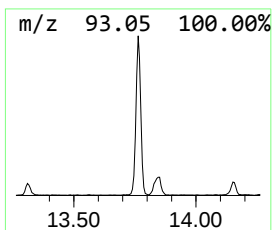
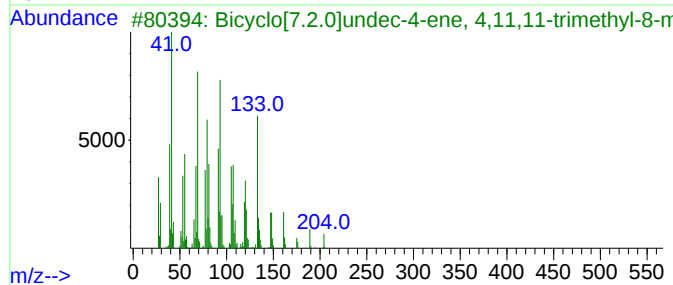
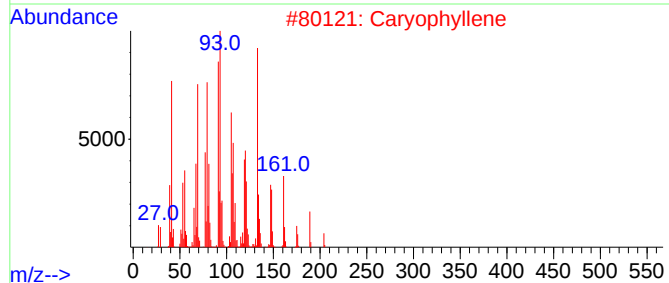
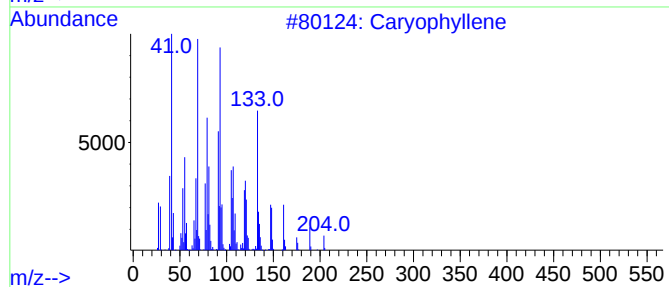
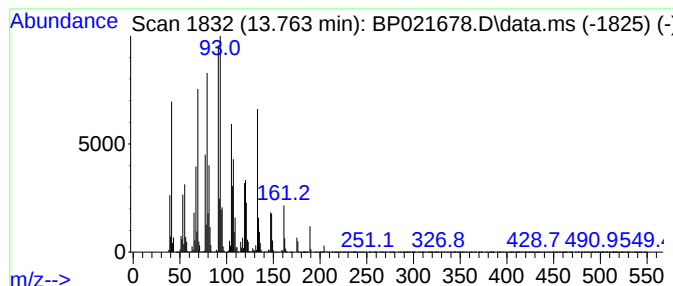
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Caryophyllene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.763	9.59 ng/ul	1833310	Acenaphthene-d10	14.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Caryophyllene	204	C15H24	000087-44-5	95
2			Caryophyllene	204	C15H24	000087-44-5	94
3			Bicyclo[7.2.0]undec-4-ene, 4,11,...	204	C15H24	000118-65-0	90
4			Caryophyllene	204	C15H24	000087-44-5	87
5			1R,3Z,9S-4,11,11-Trimethyl-8-met...	204	C15H24	1000159-39-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021678.D
Acq On : 27 Aug 2024 19:43
Operator : RC/JU
Sample : P3675-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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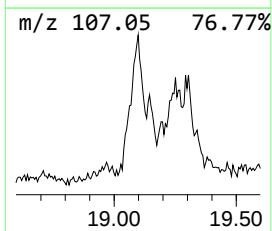
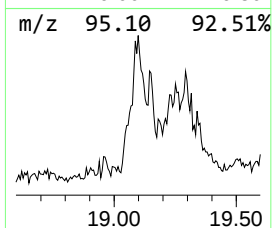
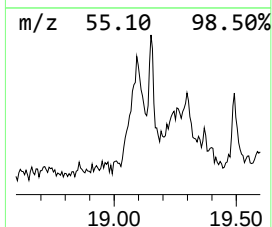
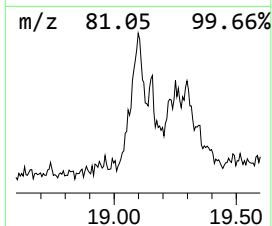
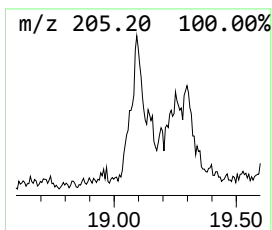
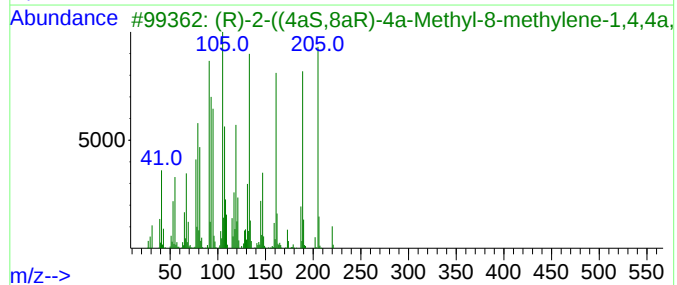
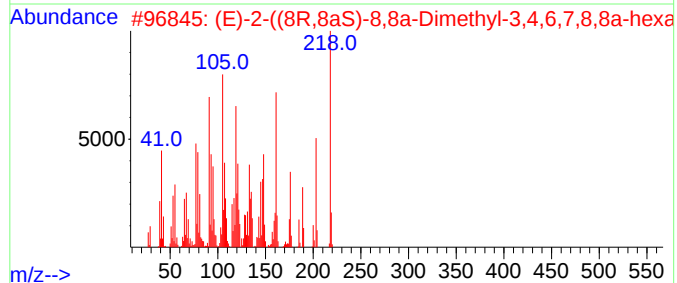
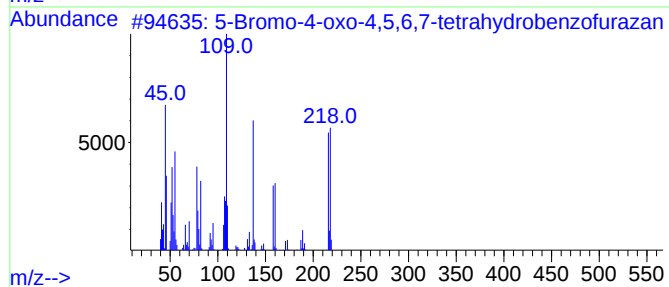
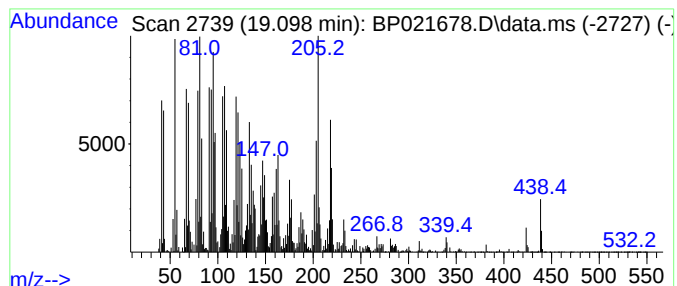
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.098	4.37 ng/ul	984585	Phenanthrene-d10	17.228

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	90
2			(E)-2-((8R,8aS)-8,8a-Dimethyl-3,...	218	C15H22O	137695-18-2	68
3			(R)-2-((4aS,8aR)-4a-Methyl-8-met...	220	C15H24O	028102-68-3	64
4			(3S,5R,8S)-3,8-Dimethyl-5-(prop-...	218	C15H22O	018374-76-0	60
5			2-Methyl-2-bornene	150	C11H18	072540-93-3	44



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021678.D
Acq On : 27 Aug 2024 19:43
Operator : RC/JU
Sample : P3675-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

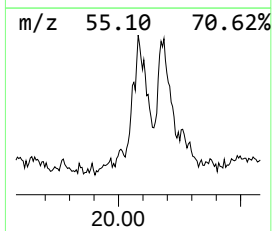
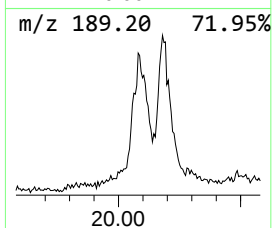
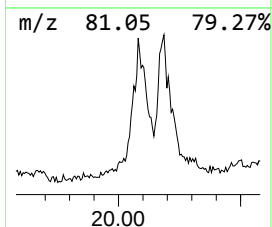
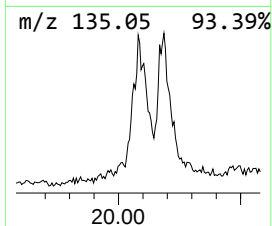
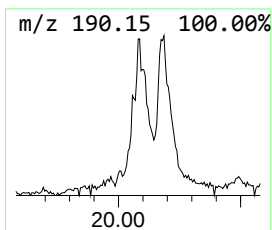
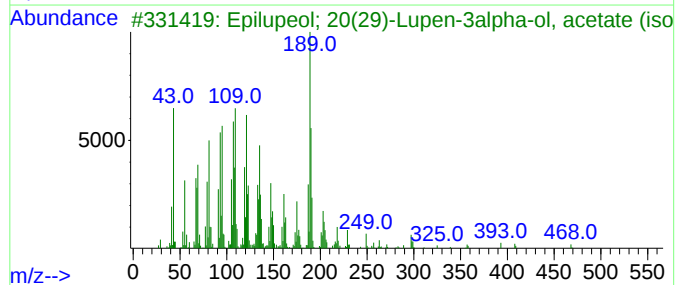
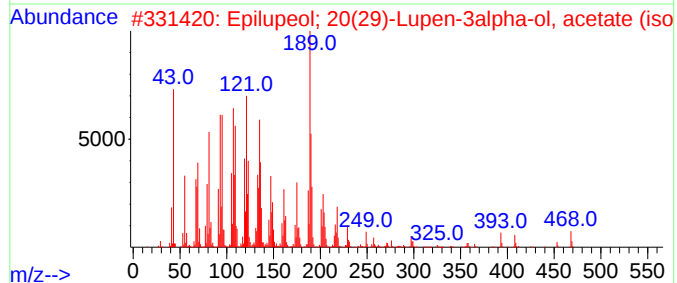
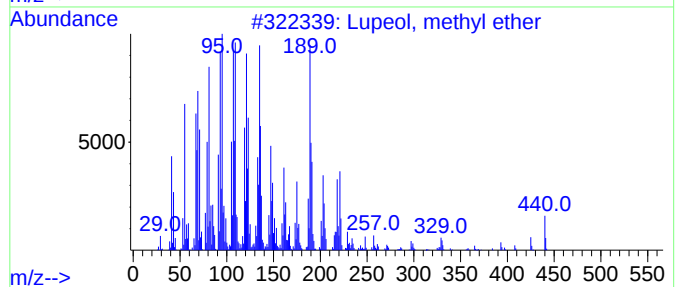
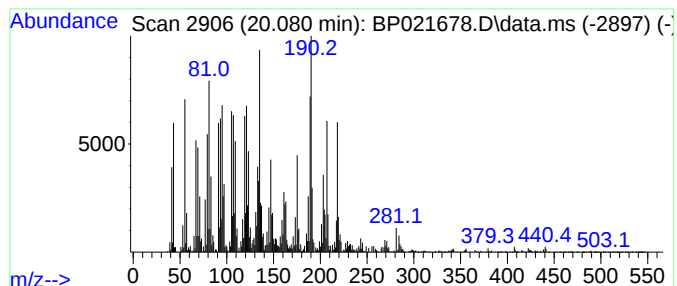
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Lupeol, methyl ether Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.080	3.58 ng/ul	931406	Chrysene-d12	21.674	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Lupeol, methyl ether	440	C31H52O	025279-38-3	78
2	Epilupeol; 20(29)-Lupen-3alpha-o...	468	C32H52O2	1000513-01-7	49
3	Epilupeol; 20(29)-Lupen-3alpha-o...	468	C32H52O2	1010513-01-8	46
4	Epilupeol; 20(29)-Lupen-3alpha-o...	426	C30H50O	1000513-01-4	45
5	4,6,6-Trimethyl-2-(3-methylbuta-...	218	C15H22O	1000190-22-2	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021678.D
Acq On : 27 Aug 2024 19:43
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Sample : P3675-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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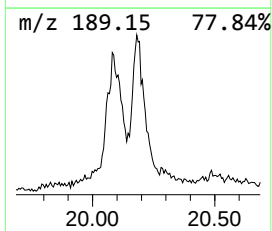
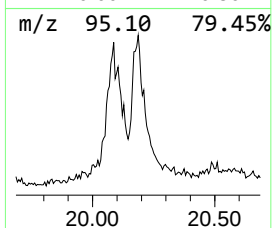
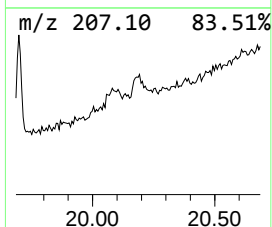
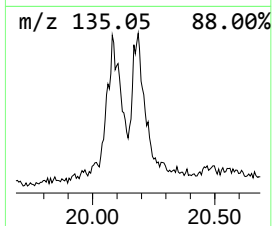
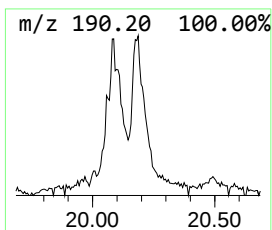
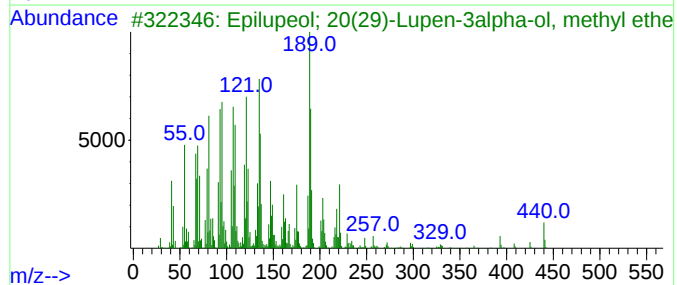
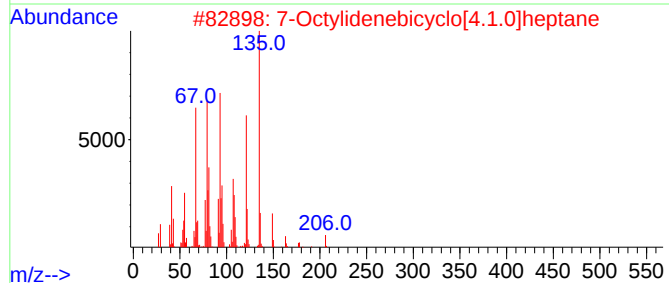
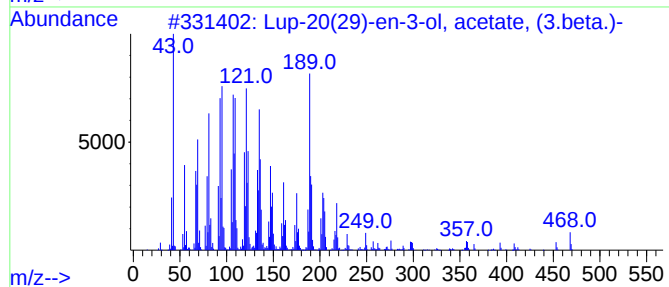
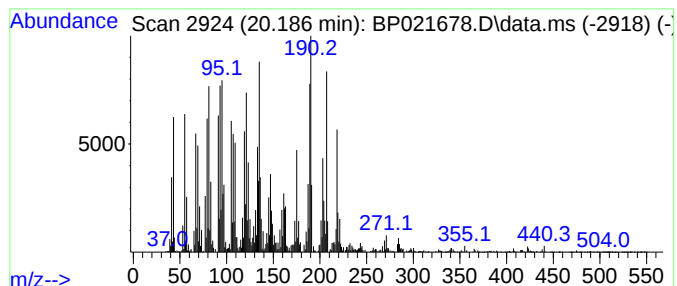
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.186	4.52 ng/ul	1174300	Chrysene-d12	21.674

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Lup-20(29)-en-3-ol, acetate, (3....	468	C32H52O2	001617-68-1	43
2			7-Octylidenebicyclo[4.1.0]heptane	206	C15H26	082253-11-0	43
3			Epilupeol; 20(29)-Lupen-3alpha-o...	440	C31H52O	1000513-01-9	42
4			7-Oxabicyclo[4.1.0]heptane, 2,2,...	218	C15H22O	070038-20-9	41
5			Boron, 1,5-cyclooctanediyl(2,4-p...	218	C13H23BN2	136704-99-9	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021678.D
Acq On : 27 Aug 2024 19:43
Operator : RC/JU
Sample : P3675-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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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
Butane, 2-metho...	3.022	49.7	ng/u1	5149800	1	7.793	2073220	20.0
Propylene Glycol	3.540	2.8	ng/u1	292438	1	7.793	2073220	20.0
Ethanol, 2-(2-e...	7.393	2.1	ng/u1	213989	1	7.793	2073220	20.0
1-Phenoxypropan...	11.322	3.1	ng/u1	443132	2	10.581	2813710	20.0
Caryophyllene	13.763	9.6	ng/u1	1833310	3	14.434	3821770	20.0
5-Bromo-4-oxo-4...	19.098	4.4	ng/u1	984585	4	17.228	4504380	20.0
Lupeol, methyl ...	20.080	3.6	ng/u1	931406	5	21.674	5198380	20.0
unknown-01	20.186	4.5	ng/u1	1174300	5	21.674	5198380	20.0