

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
 Data File : BP012747.D
 Acq On : 26 Nov 2022 02:48
 Operator : CG/JU
 Sample : N5592-05
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0AB0

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-BP112522.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012747.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.370	26	31	38	rBV	5246185	7481886	15.32%	4.139%
2	3.434	38	42	57	rVB	3154435	3959617	8.11%	2.190%
3	3.823	105	108	120	rBV	643827	908819	1.86%	0.503%
4	4.064	145	149	154	rVB	51431	71695	0.15%	0.040%
5	4.164	161	166	171	rBV	39056	49737	0.10%	0.028%
6	4.705	251	258	276	rBV	26622272	48829649	100.00%	27.011%
7	5.099	319	325	338	rBV	1830963	2528261	5.18%	1.399%
8	7.169	670	677	689	rBV	613219	965307	1.98%	0.534%
9	7.346	700	707	714	rBV	1673130	2578837	5.28%	1.427%
10	7.546	733	741	752	rBV	1691093	2660174	5.45%	1.472%
11	8.022	813	822	829	rBV	1739496	2683699	5.50%	1.485%
12	8.722	933	941	955	rBV	1614414	2563176	5.25%	1.418%
13	9.187	1012	1020	1031	rBV	1763431	2780089	5.69%	1.538%
14	9.910	1133	1143	1157	rBV	1327694	2179715	4.46%	1.206%
15	10.452	1226	1235	1251	rBV	2508317	4240142	8.68%	2.346%
16	10.840	1293	1301	1312	rBV	2402836	4064387	8.32%	2.248%
17	10.969	1315	1323	1336	rBV	1462841	2443506	5.00%	1.352%
18	12.351	1548	1558	1565	rBV	1095749	1710460	3.50%	0.946%
19	12.416	1565	1569	1575	rVB	47423	83280	0.17%	0.046%
20	14.063	1841	1849	1861	rBV	5572882	7568459	15.50%	4.187%
21	14.351	1891	1898	1917	rVB	5929808	8609479	17.63%	4.763%
22	14.657	1943	1950	1958	rBV	3918395	5790688	11.86%	3.203%
23	14.828	1973	1979	1992	rBV	423640	678804	1.39%	0.375%
24	15.063	2015	2019	2025	rBV2	68299	101209	0.21%	0.056%
25	15.651	2112	2119	2130	rBV	7992822	11556677	23.67%	6.393%
26	15.757	2130	2137	2148	rVV	1962280	2576302	5.28%	1.425%
27	15.887	2153	2159	2168	rVB2	35701	71592	0.15%	0.040%
28	17.410	2411	2418	2428	rBV	4618737	6482315	13.28%	3.586%
29	17.510	2428	2435	2448	rBV	9402685	12991840	26.61%	7.187%
30	18.281	2562	2566	2573	rBV	197283	252618	0.52%	0.140%
31	18.363	2575	2580	2585	rVV	196196	268630	0.55%	0.149%
32	19.316	2738	2742	2747	rBV	107352	142627	0.29%	0.079%
33	19.381	2749	2753	2759	rVV	99978	158700	0.33%	0.088%
34	19.510	2771	2775	2783	rBV	140537	198283	0.41%	0.110%
35	19.739	2808	2814	2832	rBV	10034712	13380114	27.40%	7.402%
36	21.498	3107	3113	3125	rBV	3304407	4685839	9.60%	2.592%

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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-BP112522.M
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37	23.851	3505	3513	3521	rBV2	4132758	8487465	17.38%	4.695%
38	24.010	3534	3540	3554	rVB2	1863034	3990410	8.17%	2.207%

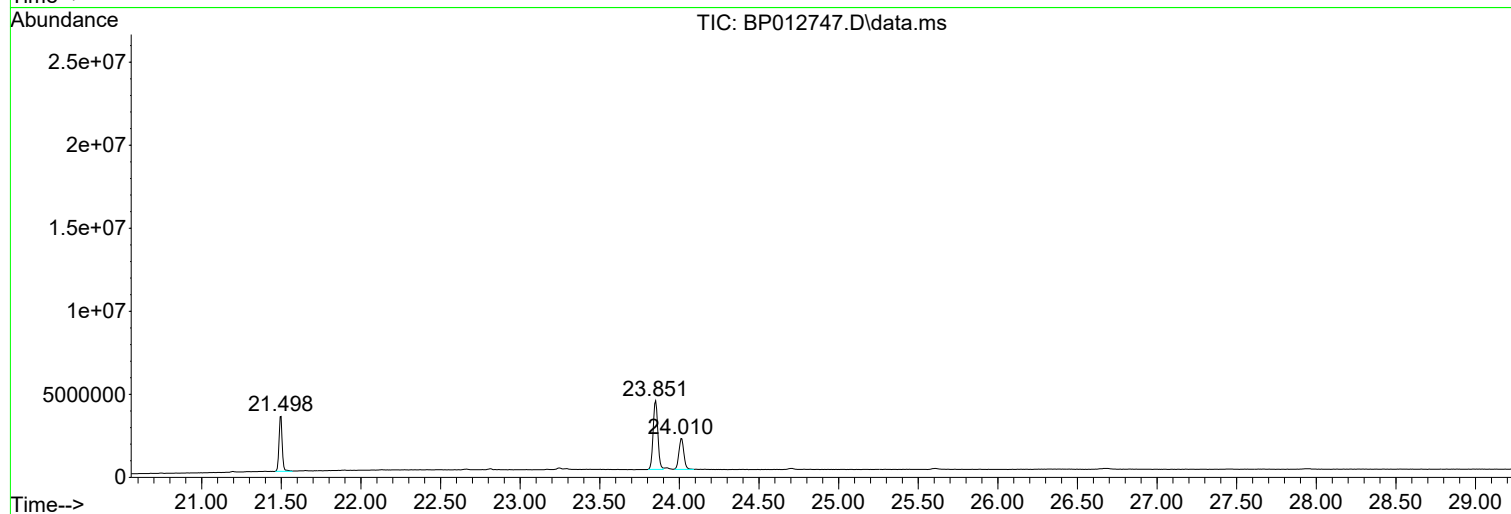
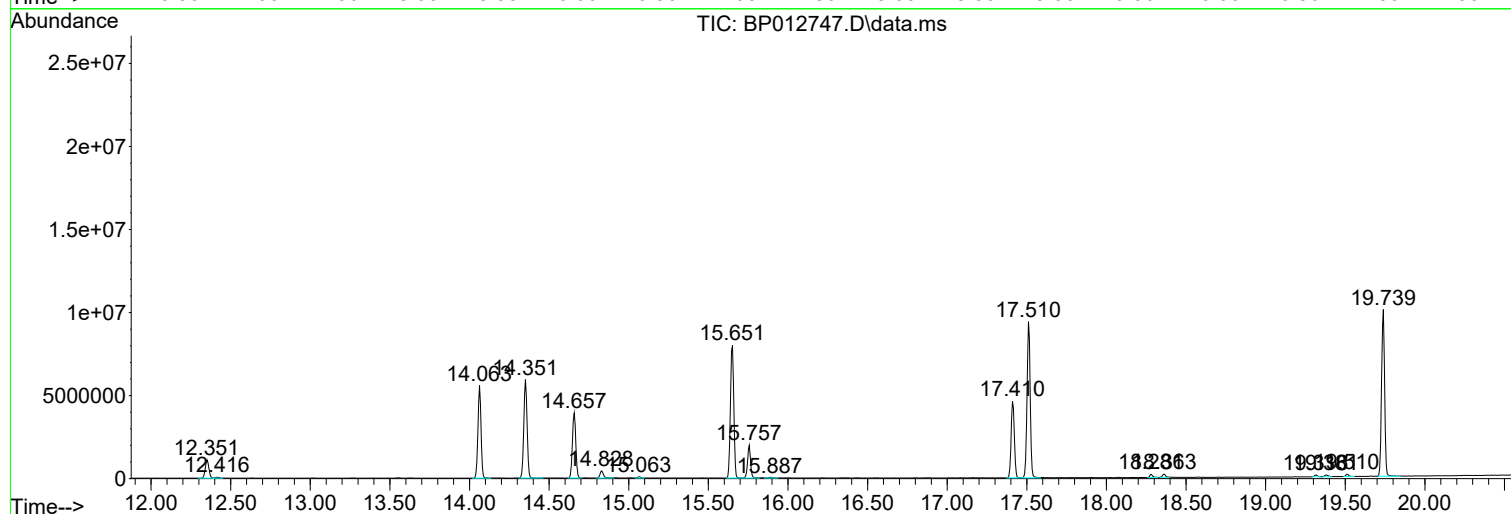
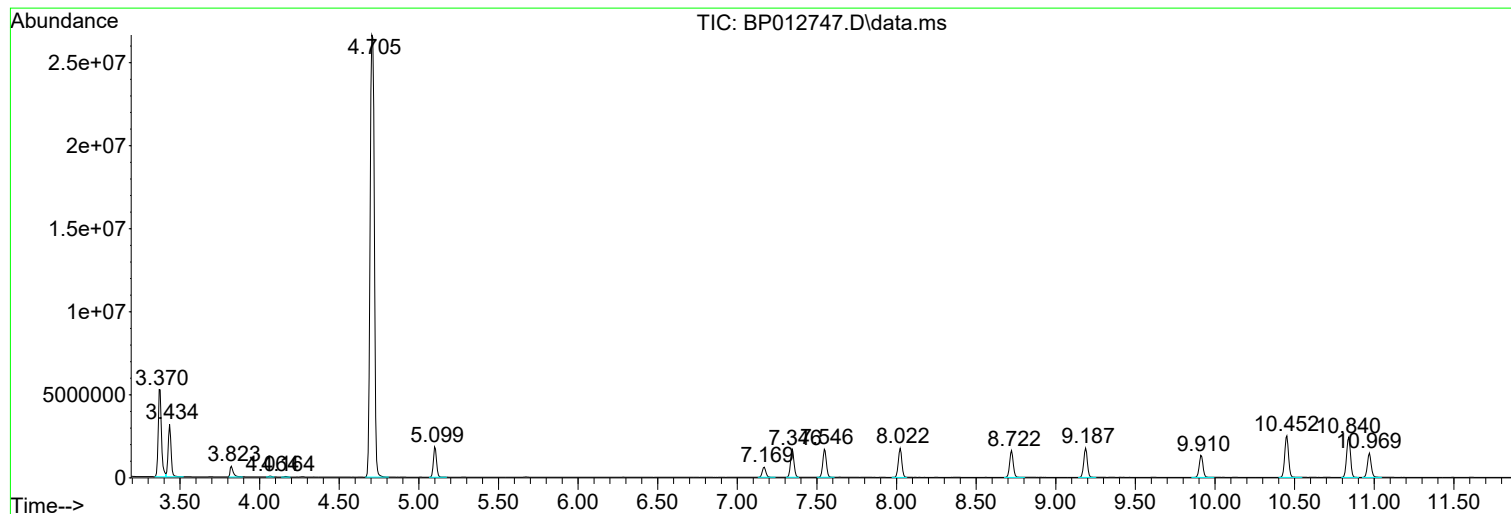
Sum of corrected areas: 180774487

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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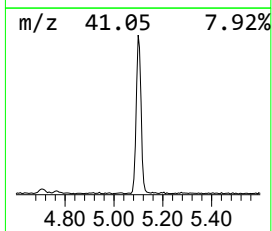
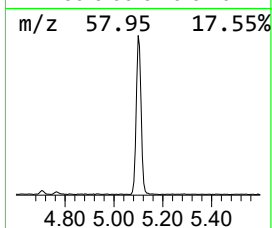
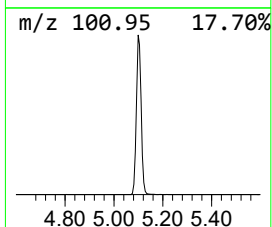
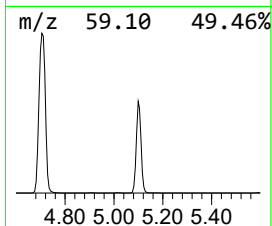
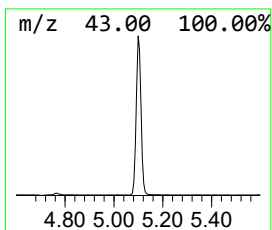
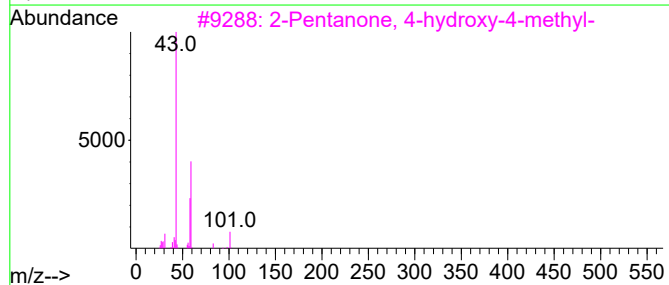
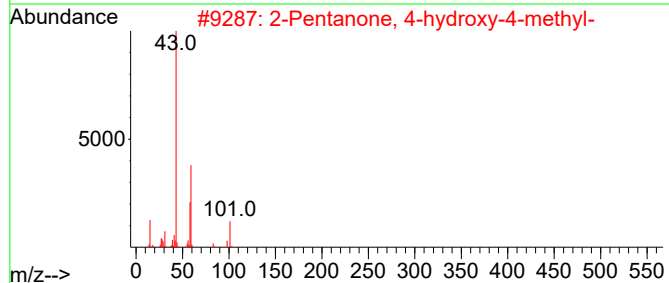
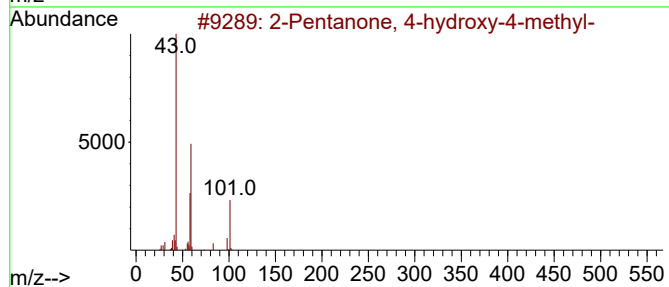
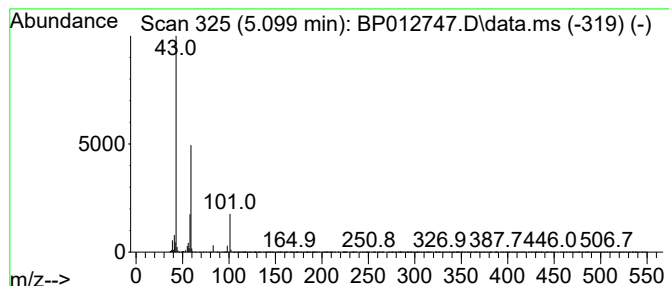
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.099	18.84 ng/ul	2528260	1,4-Dichlorobenzene-d4	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
5	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38		



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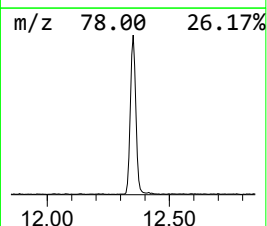
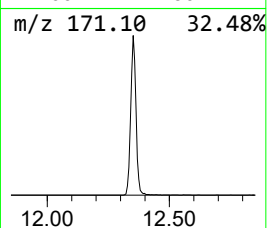
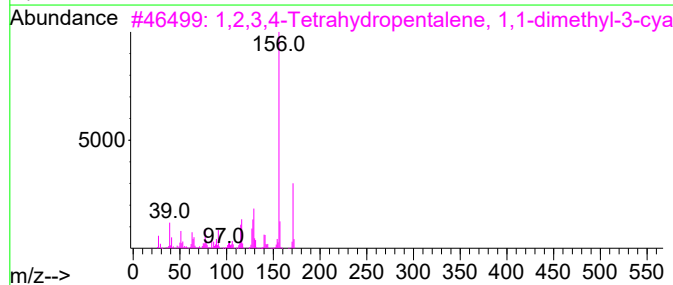
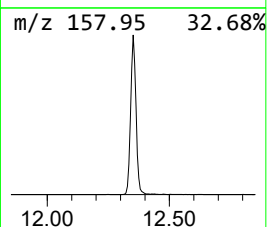
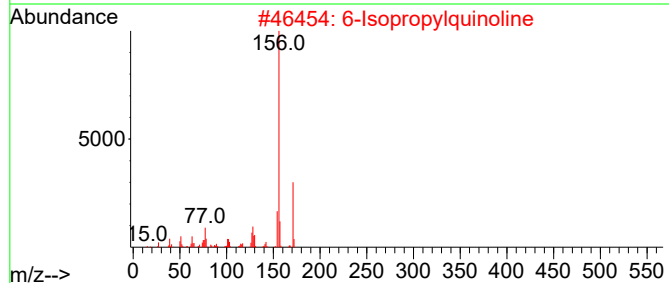
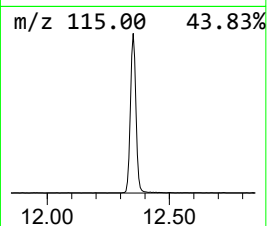
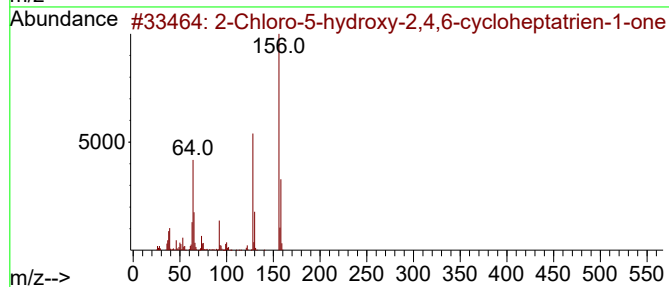
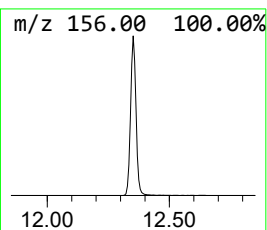
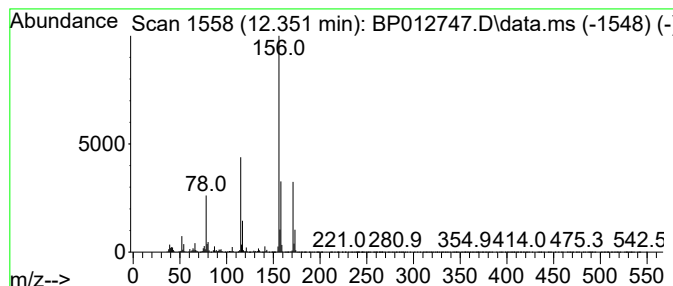
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Peak Number 3 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.352	8.42 ng/ul	1710460	Naphthalene-d8	10.840

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Chloro-5-hydroxy-2,4,6-cyclohe...	156	C7H5ClO2	007009-16-7	43		
2	6-Isopropylquinoline	171	C12H13N	000135-79-5	38		
3	1,2,3,4-Tetrahydropentalene, 1,1...	171	C12H13N	1000217-18-6	37		
4	Cyclopentanone, O-(trimethylsilyl...	171	C8H17NOSi	026208-87-7	32		
5	2-Chloro-4,6-dimethyl-3-pyridina...	156	C7H9ClN2	140413-40-7	32		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.099	18.8	ng/ul	2528260	1	8.022	2683700	20.0
unknown-01	12.352	8.4	ng/ul	1710460	2	10.840	4064390	20.0