

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
 Data File : BP017215.D
 Acq On : 18 Sep 2023 20:51
 Operator : MA/JU
 Sample : 04330-16
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZYW7

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-BP091223.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017215.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.311	34	38	49	rVB	42493	64903	1.71%	0.141%
2	3.758	109	114	121	rVB2	69251	111542	2.94%	0.242%
3	3.858	127	131	137	rVB	157667	193518	5.10%	0.420%
4	3.940	140	145	150	rVB2	90318	140952	3.71%	0.306%
5	4.058	161	165	169	rVB2	17663	25748	0.68%	0.056%
6	4.405	221	224	228	rVV3	10951	15436	0.41%	0.033%
7	4.440	228	230	237	rVB3	17643	27320	0.72%	0.059%
8	4.640	260	264	271	rVB	37127	52510	1.38%	0.114%
9	4.799	289	291	294	rVB4	5004	4970	0.13%	0.011%
10	4.987	318	323	327	rBV	52819	76289	2.01%	0.165%
11	5.028	327	330	335	rVB	50121	66469	1.75%	0.144%
12	5.111	340	344	353	rVB	22234	38220	1.01%	0.083%
13	5.452	398	402	407	rVB2	6761	8307	0.22%	0.018%
14	5.575	420	423	426	rBV	8935	10578	0.28%	0.023%
15	6.093	509	511	515	rVB5	2734	4098	0.11%	0.009%
16	6.793	626	630	633	rVB6	4003	5239	0.14%	0.011%
17	6.899	645	648	657	rVB6	8459	14002	0.37%	0.030%
18	7.099	675	682	696	rBV	666817	1118424	29.47%	2.425%
19	7.287	706	714	722	rBV	592827	896304	23.62%	1.943%
20	7.463	738	744	754	rBV	712715	1146049	30.20%	2.484%
21	7.552	754	759	765	rVV	13821	32247	0.85%	0.070%
22	7.616	768	770	777	rVB8	7422	10990	0.29%	0.024%
23	7.752	787	793	798	rVB3	7628	13439	0.35%	0.029%
24	7.946	819	826	838	rBV	1104991	1701364	44.83%	3.688%
25	8.257	876	879	883	rBV6	2330	4305	0.11%	0.009%
26	8.657	938	947	955	rBV	441345	760004	20.03%	1.648%
27	8.793	963	970	979	rVB	181240	293015	7.72%	0.635%
28	9.146	1021	1030	1039	rBV	671224	1117069	29.43%	2.422%
29	9.257	1047	1049	1053	rVB5	9903	13237	0.35%	0.029%
30	9.351	1059	1065	1071	rBV10	7001	20089	0.53%	0.044%
31	9.616	1105	1110	1113	rBV7	3308	5877	0.15%	0.013%
32	9.857	1141	1151	1163	rBV	575056	947503	24.97%	2.054%
33	9.987	1169	1173	1179	rBV2	18624	35844	0.94%	0.078%
34	10.034	1179	1181	1186	rVB6	3948	6406	0.17%	0.014%
35	10.222	1207	1213	1218	rBV5	12362	24137	0.64%	0.052%
36	10.381	1231	1240	1252	rBV	819999	1352248	35.63%	2.931%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
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37	10.657	1283	1287	1290	rVB6	4391	7286	0.19%	0.016%
38	10.769	1298	1306	1315	rBV	1500094	2460467	64.83%	5.334%
39	10.834	1315	1317	1321	rVB5	5574	6854	0.18%	0.015%
40	10.934	1327	1334	1350	rBV	441920	811440	21.38%	1.759%

41	11.322	1396	1400	1403	rBV6	5035	8745	0.23%	0.019%
42	11.528	1429	1435	1438	rBV8	7312	13637	0.36%	0.030%
43	11.587	1440	1445	1447	rBV6	5657	9238	0.24%	0.020%
44	11.798	1476	1481	1482	rBV5	5300	6699	0.18%	0.015%
45	12.169	1540	1544	1552	rVB7	6343	13822	0.36%	0.030%

46	12.351	1569	1575	1581	rVB2	17289	29422	0.78%	0.064%
47	12.922	1669	1672	1676	rBV6	3295	5225	0.14%	0.011%
48	13.057	1692	1695	1700	rVB7	5699	8370	0.22%	0.018%
49	13.175	1712	1715	1720	rVB7	8334	11495	0.30%	0.025%
50	13.292	1733	1735	1739	rBV5	2808	3891	0.10%	0.008%

51	13.357	1739	1746	1750	rBV8	8840	17342	0.46%	0.038%
52	13.445	1754	1761	1771	rBV	73661	127750	3.37%	0.277%
53	13.540	1771	1777	1778	rBV6	5301	7768	0.20%	0.017%
54	13.657	1793	1797	1799	rVB5	3743	5732	0.15%	0.012%
55	13.687	1799	1802	1806	rBV6	2642	4832	0.13%	0.010%

56	13.757	1809	1814	1815	rBV5	5094	8195	0.22%	0.018%
57	13.787	1815	1819	1824	rVB5	12191	18271	0.48%	0.040%
58	13.928	1837	1843	1846	rBV8	10147	17631	0.46%	0.038%
59	14.016	1853	1858	1868	rBV	1534112	2072240	54.60%	4.492%
60	14.222	1889	1893	1894	rBV4	3865	3837	0.10%	0.008%

61	14.304	1899	1907	1915	rBV2	1725866	2485520	65.49%	5.388%
62	14.434	1926	1929	1937	rVB5	10663	19326	0.51%	0.042%
63	14.604	1951	1958	1971	rBV	2234416	3214817	84.71%	6.969%
64	14.828	1990	1996	2009	rBV	299052	554189	14.60%	1.201%
65	14.998	2022	2025	2029	rVV4	20516	30074	0.79%	0.065%

66	15.045	2029	2033	2041	rVB3	49679	77099	2.03%	0.167%
67	15.204	2058	2060	2067	rVB8	5069	7148	0.19%	0.015%
68	15.351	2081	2085	2087	rBV3	6639	9123	0.24%	0.020%
69	15.386	2089	2091	2094	rVV3	7765	8821	0.23%	0.019%
70	15.604	2121	2128	2135	rBV	2281217	3353499	88.37%	7.270%

71	15.728	2144	2149	2155	rBV	351738	489405	12.90%	1.061%
72	15.839	2165	2168	2172	rVB5	10251	14006	0.37%	0.030%
73	16.263	2235	2240	2245	rBV2	28553	59959	1.58%	0.130%
74	16.898	2344	2348	2353	rBV4	24465	38836	1.02%	0.084%
75	17.057	2373	2375	2379	rBV2	16460	17303	0.46%	0.038%

76	17.104	2379	2383	2388	rVB4	28116	40615	1.07%	0.088%
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77	17.375	2423	2429	2440	rBV	2656820	3747448	98.75%	8.124%
78	17.475	2440	2446	2454	rVV2	2419645	3440385	90.65%	7.458%
79	17.804	2499	2502	2510	rVB3	36346	50062	1.32%	0.109%
80	17.998	2532	2535	2541	rVB2	33376	41611	1.10%	0.090%
81	18.216	2568	2572	2580	rBV2	139324	261665	6.89%	0.567%
82	19.716	2822	2827	2834	rBV	2794446	3795048	100.00%	8.227%
83	21.474	3122	3126	3133	rBV	2100014	2848877	75.07%	6.176%
84	23.833	3523	3527	3535	rVB2	1371212	2572244	67.78%	5.576%
85	24.004	3551	3556	3566	rVB	1384220	2912822	76.75%	6.315%

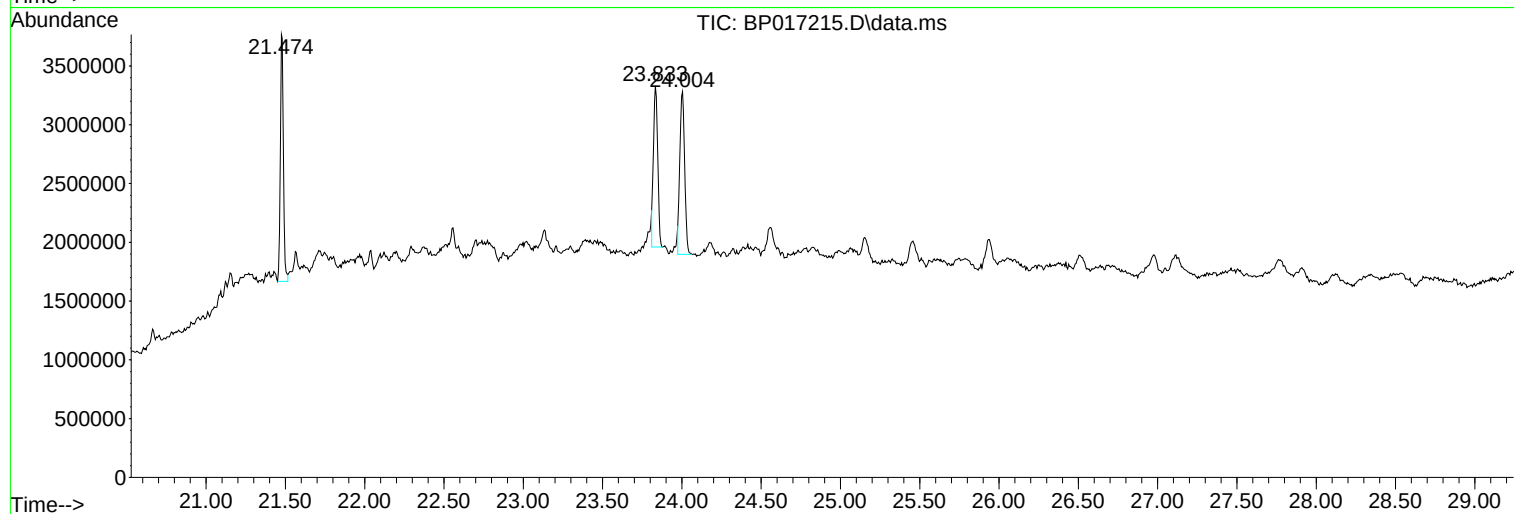
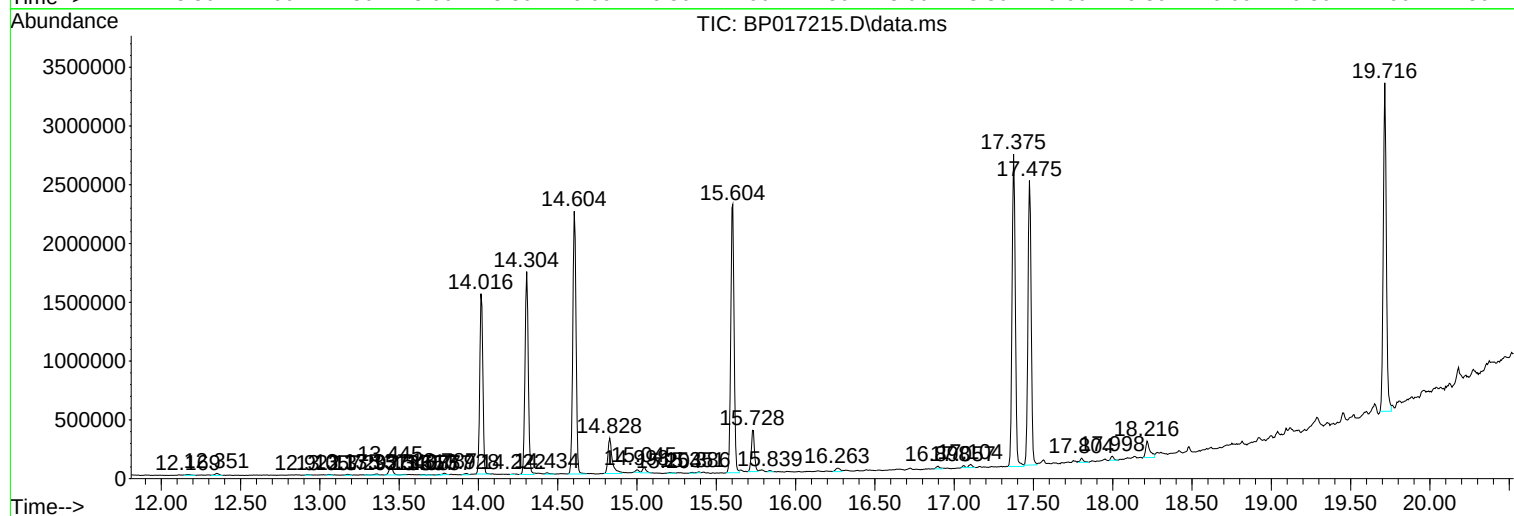
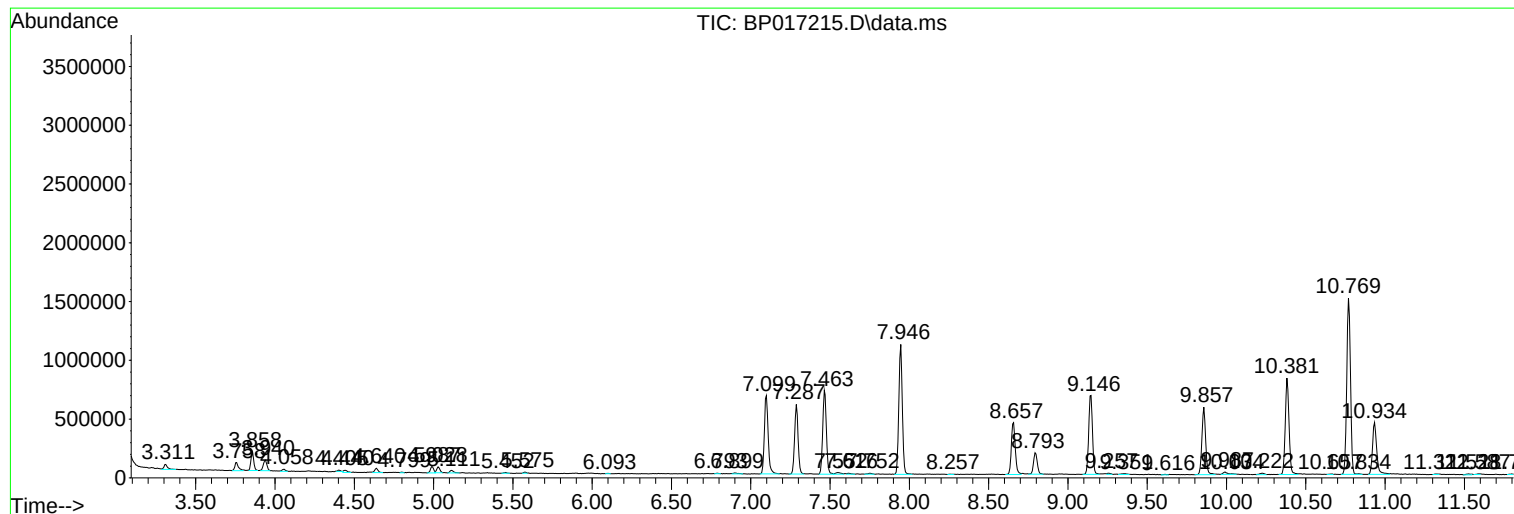
Sum of corrected areas: 46128743

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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



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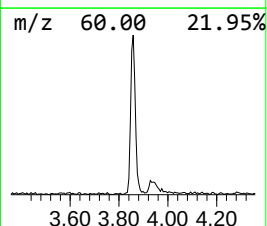
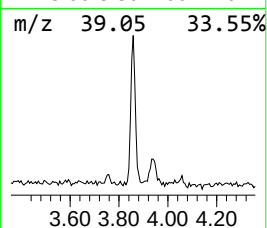
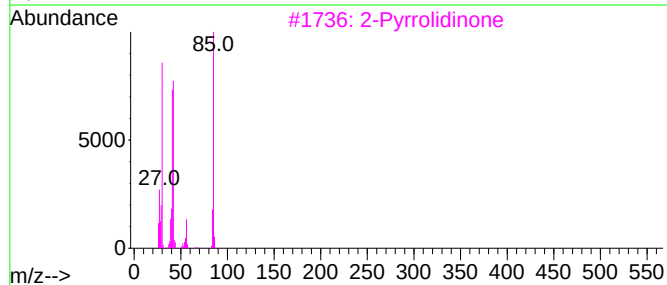
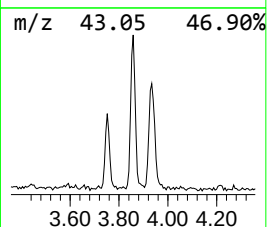
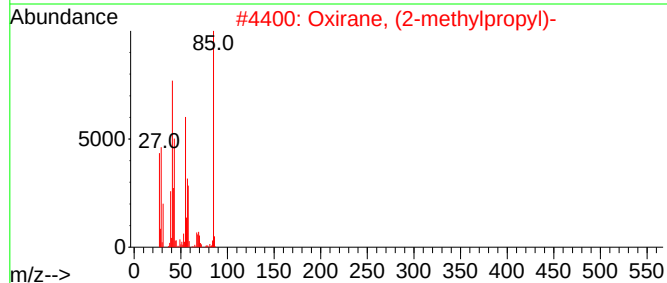
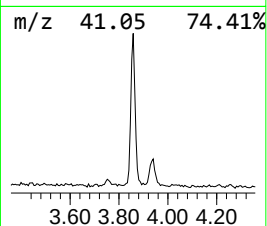
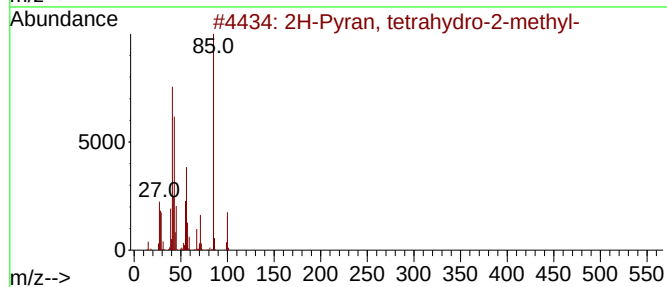
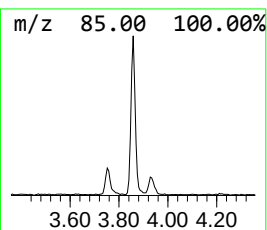
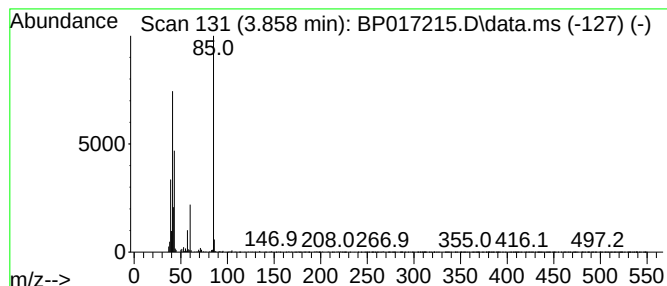
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.858	2.27 ng/ul	193518	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	38
2	Oxirane, (2-methylpropyl)-			100	C6H12O	023850-78-4	36
3	2-Pyrrolidinone			85	C4H7NO	000616-45-5	28
4	1-[(4-Fluorophenyl)sulfonyl]pipe...			244	C10H13FN2O2S	1000486-20-7	25
5	2-Pyrrolidinone			85	C4H7NO	000616-45-5	9



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	3.858	2.3	ng/u1	193518	1	7.946	1701360	20.0