

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
 Data File : BP017201.D
 Acq On : 18 Sep 2023 12:49
 Operator : MA/JU
 Sample : 04330-20
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZYL5

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: BP017201.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.310	35	38	43	rVB2	21583	26244	0.96%	0.089%
2	3.757	111	114	122	rBV	37583	51948	1.91%	0.175%
3	3.857	128	131	136	rVB	23334	28365	1.04%	0.096%
4	3.940	140	145	151	rVB	51043	77874	2.86%	0.263%
5	4.052	161	164	172	rVB	14478	20441	0.75%	0.069%
6	4.693	270	273	274	rBV3	4718	5053	0.19%	0.017%
7	4.993	319	324	329	rBV	31050	47732	1.75%	0.161%
8	5.128	342	347	353	rBV10	4992	10358	0.38%	0.035%
9	5.534	413	416	417	rBV3	2975	2765	0.10%	0.009%
10	5.581	421	424	428	rVB3	6444	8569	0.31%	0.029%
11	5.628	430	432	435	rBV4	2178	2832	0.10%	0.010%
12	5.740	448	451	454	rBV4	3805	5379	0.20%	0.018%
13	5.857	469	471	475	rBV5	3408	4084	0.15%	0.014%
14	5.899	477	478	482	rVB4	4347	4367	0.16%	0.015%
15	7.093	675	681	695	rBV	346351	567907	20.83%	1.918%
16	7.287	708	714	723	rVB	308671	463787	17.01%	1.567%
17	7.463	734	744	754	rBV	336974	538589	19.75%	1.819%
18	7.546	754	758	768	rVB	20128	39406	1.45%	0.133%
19	7.716	785	787	790	rBV3	2308	3291	0.12%	0.011%
20	7.757	790	794	802	rVB6	5744	11080	0.41%	0.037%
21	7.945	819	826	836	rBV	699257	1095856	40.19%	3.702%
22	8.287	881	884	886	rBV4	1927	2864	0.11%	0.010%
23	8.651	940	946	961	rBV	290384	473169	17.35%	1.598%
24	8.798	964	971	979	rVV	80057	134067	4.92%	0.453%
25	8.863	979	982	986	rVB6	2375	3119	0.11%	0.011%
26	8.940	994	995	998	rBV3	2939	2909	0.11%	0.010%
27	9.140	1019	1029	1037	rBV	331486	553001	20.28%	1.868%
28	9.257	1047	1049	1055	rVB7	5323	6926	0.25%	0.023%
29	9.357	1061	1066	1069	rVB5	4287	6395	0.23%	0.022%
30	9.857	1140	1151	1158	rBV	197960	333851	12.24%	1.128%
31	9.998	1170	1175	1176	rBV5	5959	9129	0.33%	0.031%
32	10.216	1209	1212	1220	rVB7	7524	13391	0.49%	0.045%
33	10.381	1231	1240	1251	rBV	336184	572494	21.00%	1.934%
34	10.769	1297	1306	1313	rBV	965080	1599299	58.65%	5.402%
35	10.822	1313	1315	1321	rVB5	8433	11333	0.42%	0.038%
36	10.934	1326	1334	1345	rBV	363333	643885	23.61%	2.175%

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

37	11.316	1395	1399	1400	rVV4	5108	5653	0.21%	0.019%
38	11.334	1400	1402	1406	rVV5	4558	5862	0.21%	0.020%
39	11.398	1409	1413	1417	rVB7	3042	4129	0.15%	0.014%
40	11.604	1444	1448	1450	rVB5	2816	3231	0.12%	0.011%

41	11.810	1480	1483	1485	rBV4	1955	2927	0.11%	0.010%
42	12.116	1532	1535	1539	rVB6	4098	5076	0.19%	0.017%
43	12.181	1543	1546	1548	rBV4	3318	3482	0.13%	0.012%
44	12.351	1571	1575	1579	rVB5	9043	12687	0.47%	0.043%
45	12.651	1623	1626	1629	rBV5	2797	4024	0.15%	0.014%

46	12.869	1661	1663	1666	rVB4	3370	2925	0.11%	0.010%
47	12.904	1666	1669	1671	rBV3	2475	2859	0.10%	0.010%
48	12.933	1671	1674	1677	rVB4	2952	3692	0.14%	0.012%
49	13.280	1729	1733	1736	rBV6	2854	3739	0.14%	0.013%
50	13.357	1739	1746	1749	rBV5	10926	13323	0.49%	0.045%

51	13.451	1754	1762	1772	rBV	674345	1062397	38.96%	3.588%
52	13.645	1792	1795	1800	rVB6	2529	3192	0.12%	0.011%
53	13.780	1815	1818	1824	rBV8	4706	7155	0.26%	0.024%
54	13.922	1837	1842	1847	rBV9	5409	8905	0.33%	0.030%
55	13.963	1847	1849	1853	rBV4	2950	4031	0.15%	0.014%

56	14.016	1853	1858	1870	rBV	824023	1118616	41.02%	3.778%
57	14.145	1877	1880	1885	rVB7	3378	5871	0.22%	0.020%
58	14.304	1900	1907	1919	rBV	901021	1280822	46.97%	4.326%
59	14.433	1926	1929	1936	rVB7	9420	17148	0.63%	0.058%
60	14.604	1951	1958	1968	rBV	1530876	2154537	79.01%	7.277%

61	14.845	1996	1999	2002	rBV4	4479	6861	0.25%	0.023%
62	15.004	2023	2026	2029	rBV5	2789	3928	0.14%	0.013%
63	15.051	2030	2034	2041	rVB9	9331	15217	0.56%	0.051%
64	15.210	2058	2061	2066	rVB7	2448	3294	0.12%	0.011%
65	15.269	2069	2071	2075	rVB5	5953	5949	0.22%	0.020%

66	15.339	2081	2083	2087	rBV5	4751	6266	0.23%	0.021%
67	15.410	2088	2095	2099	rVV2	61308	100378	3.68%	0.339%
68	15.445	2099	2101	2108	rVB3	19480	25497	0.94%	0.086%
69	15.504	2108	2111	2117	rVB2	16226	20084	0.74%	0.068%
70	15.598	2120	2127	2144	rBV	1191371	1795212	65.84%	6.064%

71	15.792	2155	2160	2164	rVB8	6934	12438	0.46%	0.042%
72	15.839	2166	2168	2171	rVB4	4309	5086	0.19%	0.017%
73	16.027	2196	2200	2204	rBV7	3002	4386	0.16%	0.015%
74	16.204	2226	2230	2235	rVB3	20487	30561	1.12%	0.103%
75	16.263	2235	2240	2246	rBV3	18521	32624	1.20%	0.110%

76	16.610	2294	2299	2304	rBV2	16095	27003	0.99%	0.091%
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 Title : SVOA CALIBRATION

77	16.721	2314	2318	2321	rBV6	9976	14610	0.54%	0.049%
78	16.798	2327	2331	2334	rBV5	9717	11759	0.43%	0.040%
79	16.898	2341	2348	2350	rBV2	29355	54721	2.01%	0.185%
80	16.927	2350	2353	2362	rVB	62779	102820	3.77%	0.347%
81	17.063	2372	2376	2379	rVB2	12260	17328	0.64%	0.059%
82	17.374	2423	2429	2435	rBV	1882464	2633715	96.59%	8.896%
83	17.474	2440	2446	2456	rVV	1374453	1900273	69.69%	6.419%
84	17.563	2458	2461	2467	rVB2	52282	74508	2.73%	0.252%
85	17.804	2499	2502	2508	rVB3	21295	24469	0.90%	0.083%
86	17.998	2531	2535	2541	rBV3	27781	40601	1.49%	0.137%
87	18.216	2568	2572	2580	rBV8	42058	68187	2.50%	0.230%
88	18.480	2614	2617	2621	rVB3	23183	25484	0.93%	0.086%
89	19.104	2719	2723	2725	rBV4	34696	51393	1.88%	0.174%
90	19.133	2725	2728	2732	rVV2	144743	163020	5.98%	0.551%
91	19.351	2763	2765	2768	rBV3	30678	33665	1.23%	0.114%
92	19.715	2821	2827	2838	rBV	2020500	2726756	100.00%	9.210%
93	21.474	3121	3126	3131	rBV	1827669	2403902	88.16%	8.120%
94	23.827	3521	3526	3542	rVB2	922594	1961513	71.94%	6.625%
95	23.998	3549	3555	3565	rVB	1008788	2084171	76.43%	7.040%

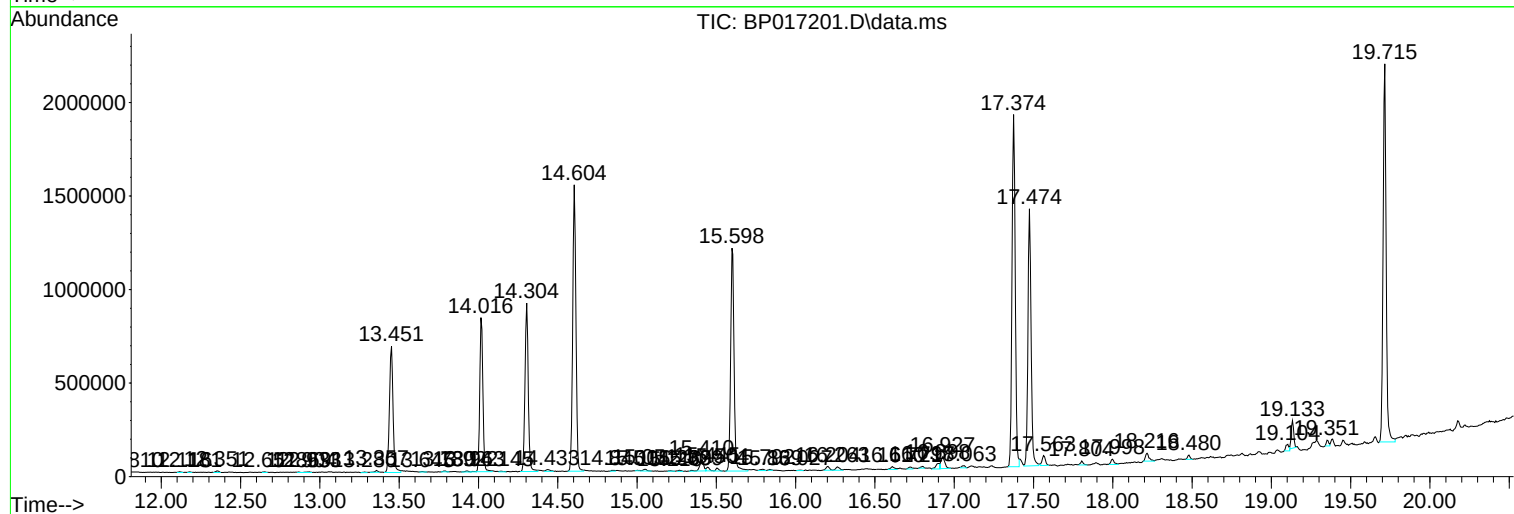
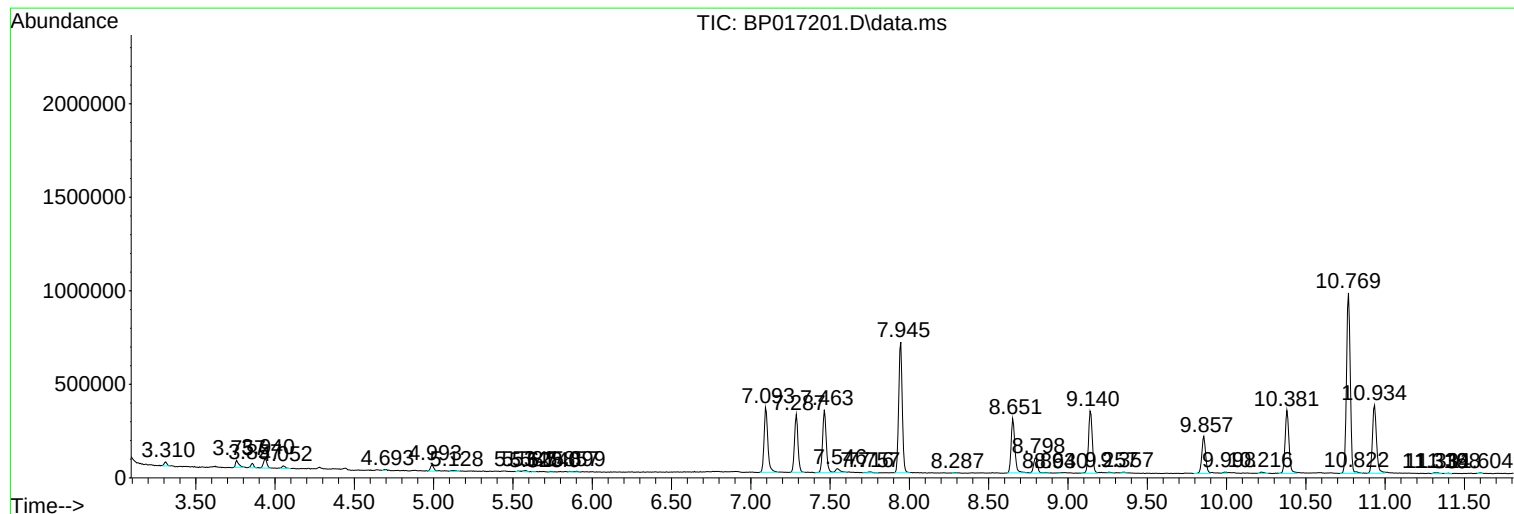
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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



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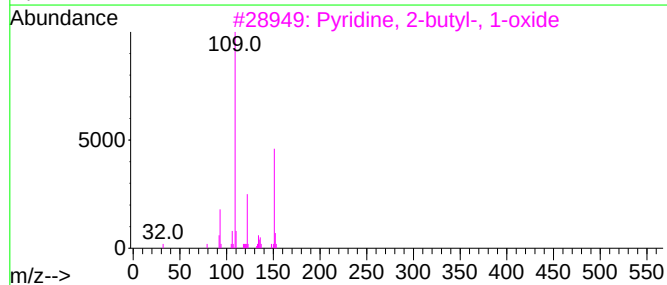
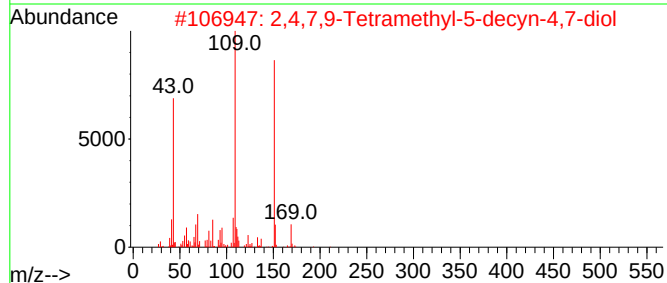
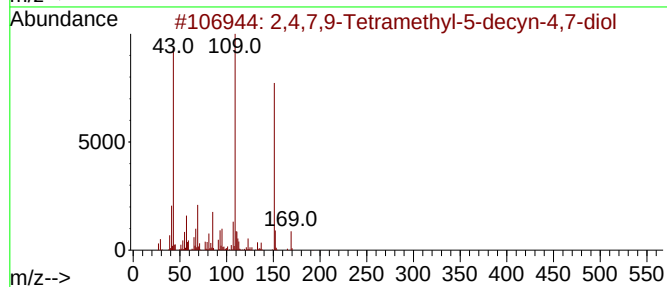
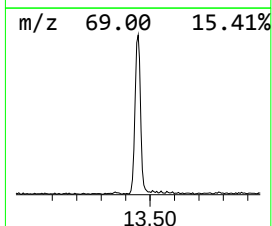
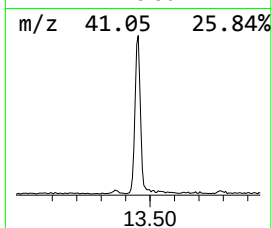
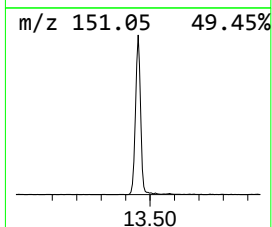
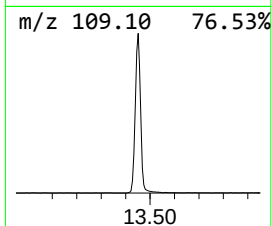
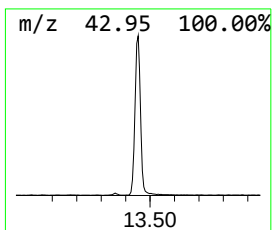
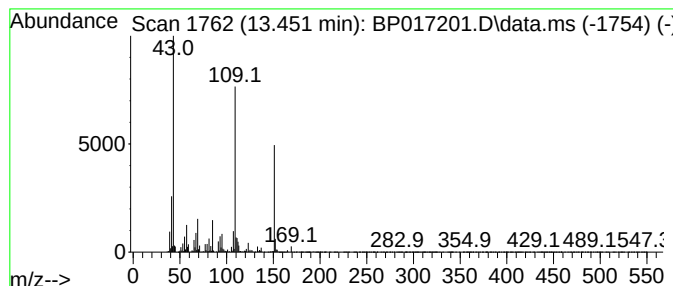
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.451	9.86 ng/ul	1062400	Acenaphthene-d10		14.604	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...		226	C14H26O2	000126-86-3	91
2	2,4,7,9-Tetramethyl-5-decyn-4,7-...		226	C14H26O2	000126-86-3	90
3	Pyridine, 2-butyl-, 1-oxide		151	C9H13NO	031396-32-4	59
4	3-Pyridinol, 4-methyl-, acetate ...		151	C8H9NO2	001006-96-8	59
5	Acetaminophen		151	C8H9NO2	000103-90-2	59



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2,4,7,9-Tetrame...	13.451	9.9	ng/ul	1062400	3	14.604	2154540	20.0