

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
 Data File : BP023509.D
 Acq On : 18 Dec 2024 22:24
 Operator : RC/JU
 Sample : P5235-10 10X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AW0

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-BP112624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023509.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.093	15	18	21	rBV2	49380	63642	0.82%	0.087%
2	3.128	21	24	34	rVB	357638	438786	5.64%	0.602%
3	3.428	72	75	80	rVB2	36063	51054	0.66%	0.070%
4	3.522	88	91	93	rBV	165980	236492	3.04%	0.324%
5	3.793	133	137	148	rVB2	62018	110483	1.42%	0.152%
6	4.064	179	183	189	rVB	95848	125524	1.61%	0.172%
7	4.140	192	196	201	rBV7	13302	22086	0.28%	0.030%
8	4.246	209	214	220	rBV	94694	135123	1.74%	0.185%
9	4.311	221	225	232	rVB	48969	74591	0.96%	0.102%
10	4.681	282	288	293	rBV	174933	244518	3.14%	0.335%
11	4.728	293	296	304	rVB	33852	57259	0.74%	0.079%
12	5.269	385	388	395	rVB3	11134	17826	0.23%	0.024%
13	5.528	428	432	436	rBV5	6860	10354	0.13%	0.014%
14	5.575	438	440	446	rVV7	5378	8041	0.10%	0.011%
15	5.628	446	449	457	rVV7	6848	11841	0.15%	0.016%
16	6.434	581	586	593	rVB	28638	48205	0.62%	0.066%
17	6.669	623	626	634	rVB10	5515	9528	0.12%	0.013%
18	6.752	634	640	656	rBV	148729	378879	4.87%	0.520%
19	6.881	656	662	677	rVV	873211	1419530	18.23%	1.947%
20	6.999	679	682	689	rVB5	8299	15587	0.20%	0.021%
21	7.081	689	696	711	rBV	778118	1427300	18.33%	1.958%
22	7.299	729	733	739	rBV3	11696	18768	0.24%	0.026%
23	7.528	763	772	780	rBV	839343	1384513	17.78%	1.899%
24	7.593	781	783	800	rVB3	20395	57573	0.74%	0.079%
25	7.863	824	829	835	rBV5	12866	25655	0.33%	0.035%
26	8.252	888	895	910	rBV	592433	1241615	15.95%	1.703%
27	8.363	911	914	918	rVV2	24767	38206	0.49%	0.052%
28	8.428	921	925	934	rVB7	11151	22816	0.29%	0.031%
29	8.675	958	967	981	rBV	885488	1713870	22.01%	2.351%
30	8.834	990	994	997	rBV5	4958	8247	0.11%	0.011%
31	8.881	997	1002	1006	rVV2	29077	51042	0.66%	0.070%
32	8.916	1006	1008	1014	rVV3	13196	23511	0.30%	0.032%
33	8.975	1014	1018	1025	rVB4	7446	17576	0.23%	0.024%
34	9.110	1035	1041	1046	rVB4	11174	17463	0.22%	0.024%
35	9.240	1059	1063	1070	rVB9	8619	19164	0.25%	0.026%
36	9.381	1079	1087	1104	rBV	799802	1473878	18.93%	2.022%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
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37	9.699	1133	1141	1148	rBV8	17773	42413	0.54%	0.058%
38	9.816	1157	1161	1172	rVB8	5569	13255	0.17%	0.018%
39	9.934	1172	1181	1208	rBV2	1392522	3193003	41.01%	4.380%
40	10.216	1224	1229	1232	rBV2	84933	128034	1.64%	0.176%
41	10.263	1232	1237	1253	rVB	1173122	1917422	24.63%	2.630%
42	10.428	1257	1265	1282	rBV	695309	1586561	20.38%	2.176%
43	10.546	1282	1285	1296	rVB6	20146	41297	0.53%	0.057%
44	10.804	1323	1329	1334	rBV	199086	349595	4.49%	0.480%
45	10.863	1334	1339	1351	rVB	268396	517262	6.64%	0.710%
46	11.104	1374	1380	1396	rBV2	146074	279250	3.59%	0.383%
47	11.481	1439	1444	1449	rBV2	18552	27439	0.35%	0.038%
48	11.598	1459	1464	1468	rBV8	6764	10785	0.14%	0.015%
49	11.657	1468	1474	1477	rBV6	7651	14916	0.19%	0.020%
50	11.869	1506	1510	1518	rVB3	15822	30468	0.39%	0.042%
51	11.981	1521	1529	1538	rBV4	22391	62309	0.80%	0.085%
52	12.416	1599	1603	1608	rBV7	9718	15795	0.20%	0.022%
53	12.681	1637	1648	1661	rBV4	66360	152913	1.96%	0.210%
54	12.969	1691	1697	1700	rBV3	18750	33376	0.43%	0.046%
55	13.022	1700	1706	1716	rVB	155194	292926	3.76%	0.402%
56	13.145	1721	1727	1738	rVV2	37088	87413	1.12%	0.120%
57	13.245	1738	1744	1747	rVV6	19544	34571	0.44%	0.047%
58	13.275	1747	1749	1760	rVB10	11577	26428	0.34%	0.036%
59	13.398	1766	1770	1774	rVB5	13318	18152	0.23%	0.025%
60	13.557	1790	1797	1810	rVV	2508435	3615512	46.44%	4.959%
61	13.828	1831	1843	1860	rBV	2499951	4019832	51.63%	5.514%
62	13.992	1864	1871	1883	rVV2	36269	92141	1.18%	0.126%
63	14.140	1889	1896	1907	rBV	1756525	2818768	36.20%	3.866%
64	14.287	1917	1921	1927	rVB7	8273	15490	0.20%	0.021%
65	14.492	1950	1956	1965	rBV3	65291	210238	2.70%	0.288%
66	14.957	2030	2035	2038	rBV7	11264	15758	0.20%	0.022%
67	15.022	2040	2046	2054	rBV2	33827	73762	0.95%	0.101%
68	15.151	2062	2068	2086	rVB	3681889	5337521	68.55%	7.321%
69	15.322	2091	2097	2108	rBV	1005490	1701027	21.85%	2.333%
70	15.539	2129	2134	2141	rBV	72406	117278	1.51%	0.161%
71	16.475	2289	2293	2294	rBV2	9043	11769	0.15%	0.016%
72	16.739	2334	2338	2343	rVB2	11113	17037	0.22%	0.023%
73	16.951	2368	2374	2385	rBV	2341601	3611509	46.39%	4.954%
74	17.057	2385	2392	2413	rVV	3884216	6249635	80.27%	8.573%
75	17.210	2415	2418	2423	rVB5	8068	12761	0.16%	0.018%
76	17.639	2484	2491	2497	rVB7	12616	29442	0.38%	0.040%

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 Title : SVOA CALIBRATION

77	17.710	2501	2503	2506	rBV2	9640	12255	0.16%	0.017%
78	17.810	2517	2520	2527	rBV5	12501	21518	0.28%	0.030%
79	17.881	2528	2532	2540	rVV	171537	299918	3.85%	0.411%
80	17.945	2541	2543	2550	rVB	31703	42923	0.55%	0.059%
81	18.069	2558	2564	2565	rBV6	7219	13879	0.18%	0.019%
82	18.275	2596	2599	2604	rVB	18349	26481	0.34%	0.036%
83	18.975	2714	2718	2724	rBV	75550	97336	1.25%	0.134%
84	19.057	2727	2732	2736	rVV	53252	91347	1.17%	0.125%
85	19.092	2736	2738	2747	rVB10	28918	56699	0.73%	0.078%
86	19.222	2756	2760	2770	rBV2	121384	229509	2.95%	0.315%
87	19.433	2789	2796	2805	rBV2	5123243	7785780	100.00%	10.680%
88	19.510	2806	2809	2816	rVB	52067	98876	1.27%	0.136%
89	21.380	3121	3127	3139	rBV	2746972	4207761	54.04%	5.772%
90	24.333	3620	3629	3646	rVB	2990933	7683059	98.68%	10.539%
91	24.545	3657	3665	3684	rVB	1549141	4389575	56.38%	6.021%

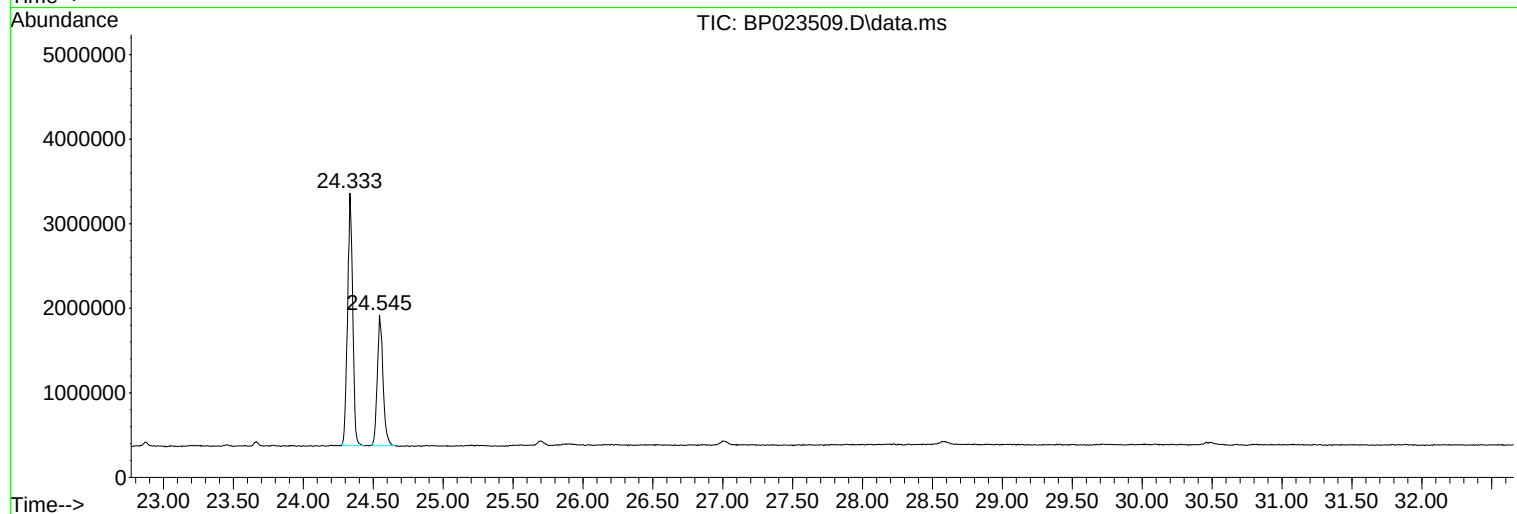
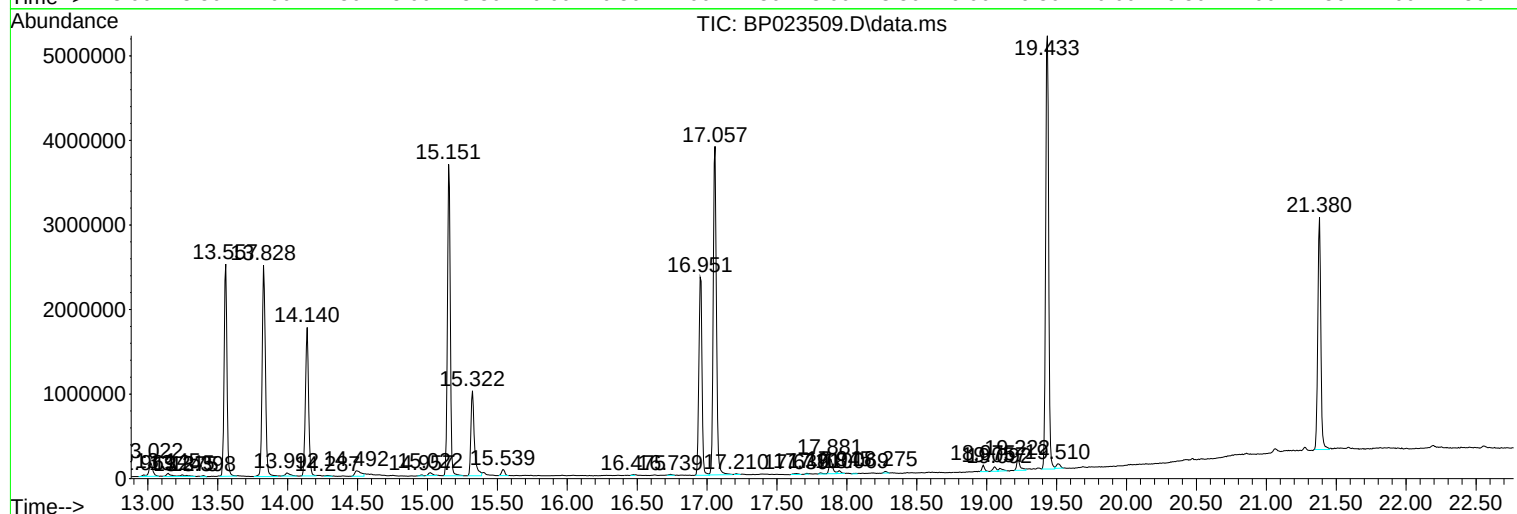
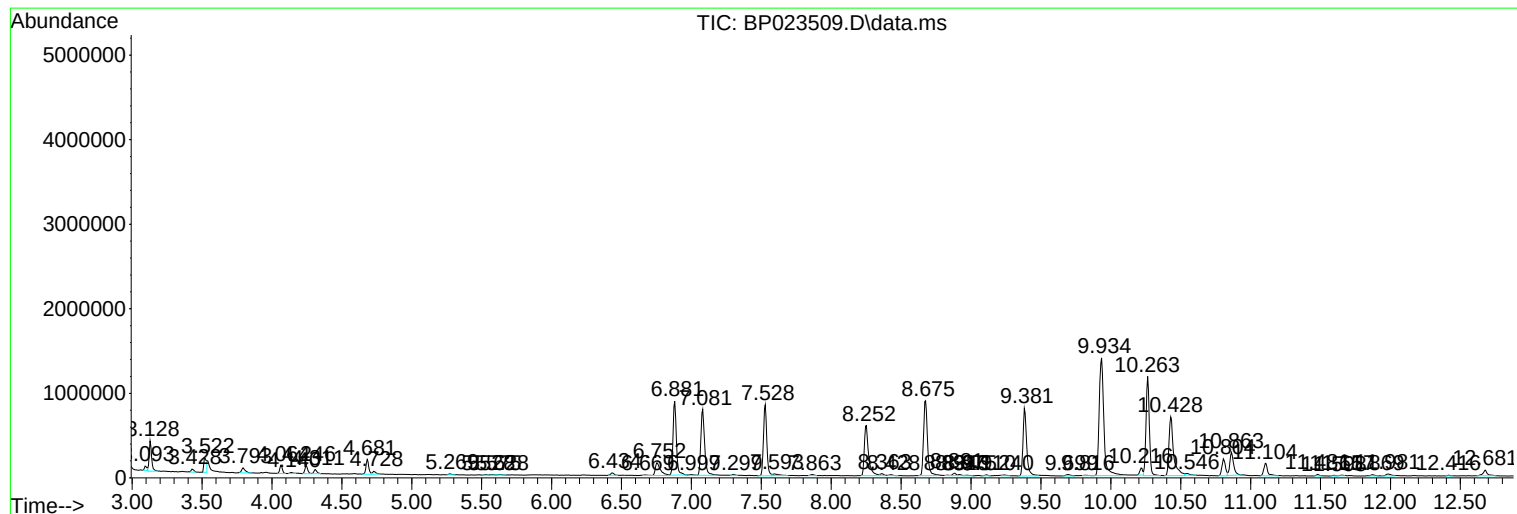
Sum of corrected areas: 72902525

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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



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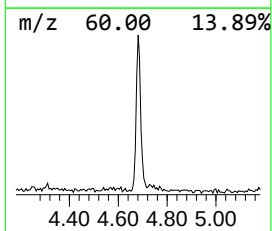
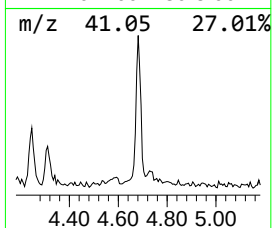
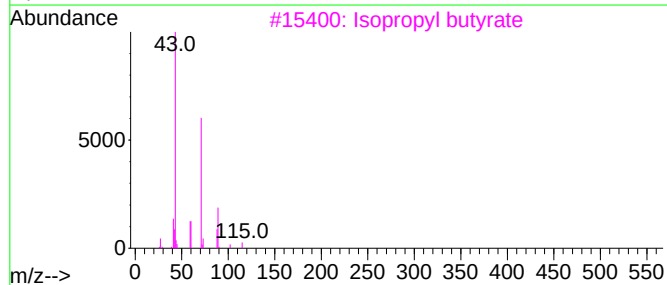
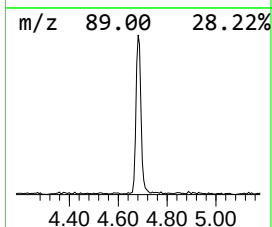
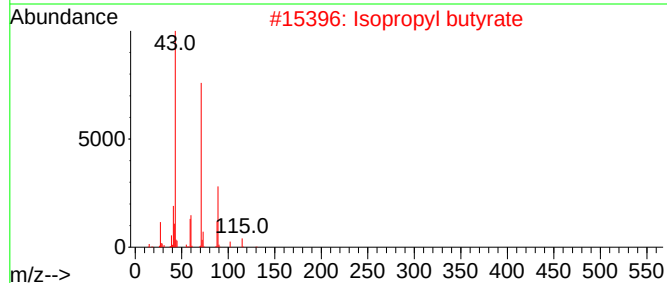
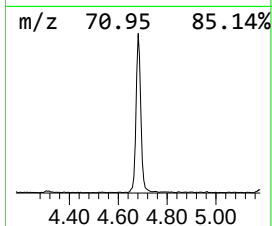
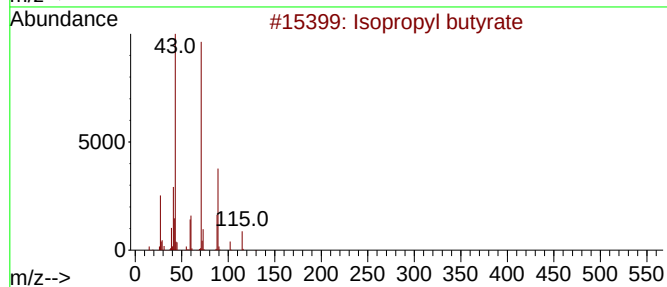
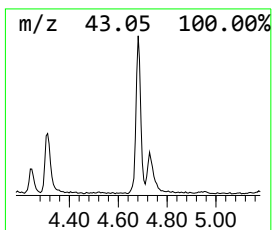
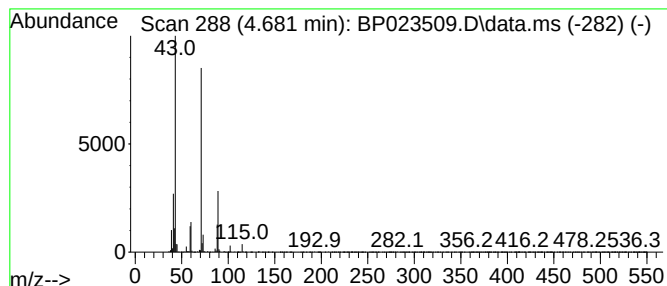
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Isopropyl butyrate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.681	3.53 ng/ul	244518	1,4-Dichlorobenzene-d4	7.528

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl butyrate	130	C7H14O2	000638-11-9	90
2			Isopropyl butyrate	130	C7H14O2	000638-11-9	90
3			Isopropyl butyrate	130	C7H14O2	000638-11-9	78
4			Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	64
5			Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	59



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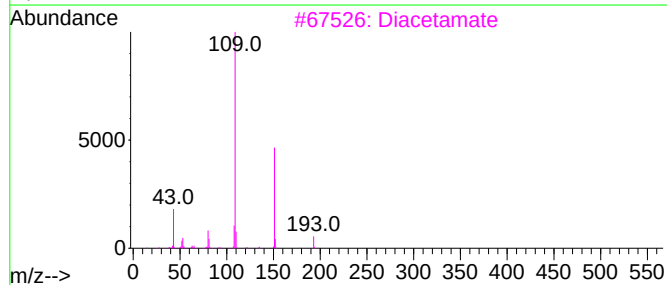
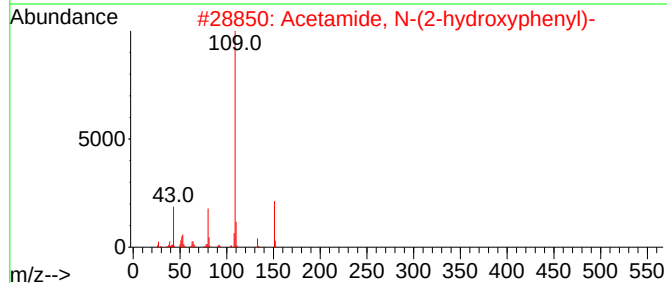
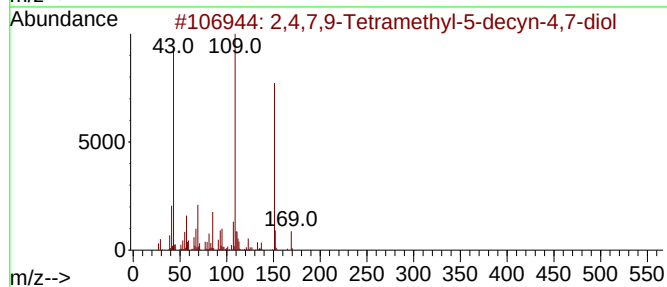
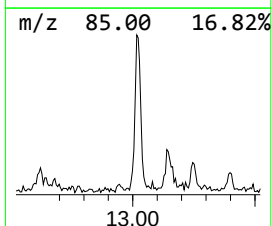
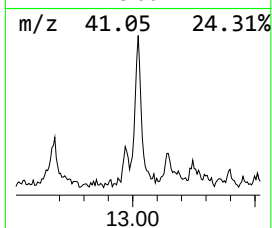
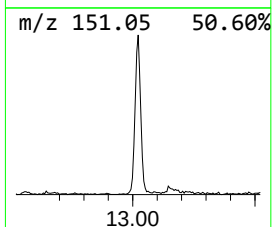
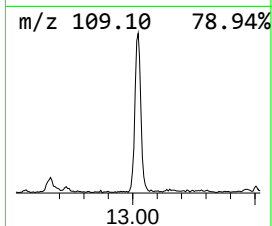
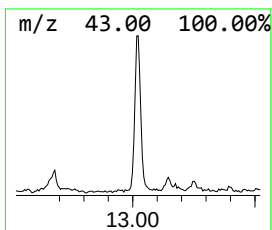
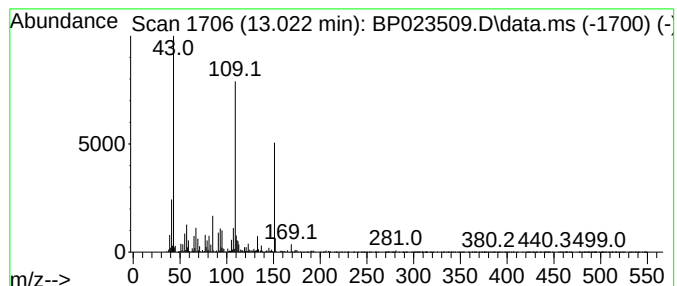
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.022	2.08 ng/ul	292926	Acenaphthene-d10	14.140	
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	Acetamide, N-(2-hydroxyphenyl)-	151	C8H9NO2	000614-80-2	59
3	Diacetamate	193	C10H11NO3	002623-33-8	53
4	Metacetamol	151	C8H9NO2	000621-42-1	53
5	m-Isopropoxyaniline	151	C9H13NO	041406-00-2	53



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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
Isopropyl butyrate	4.681	3.5	ng/u1	244518	1	7.528	1384510	20.0
2,4,7,9-Tetrame...	13.022	2.1	ng/u1	292926	3	14.140	2818770	20.0