

mohammad
7/27/2021 12:05:57 PM

[illegible]

Quantitation Report (Qedit)

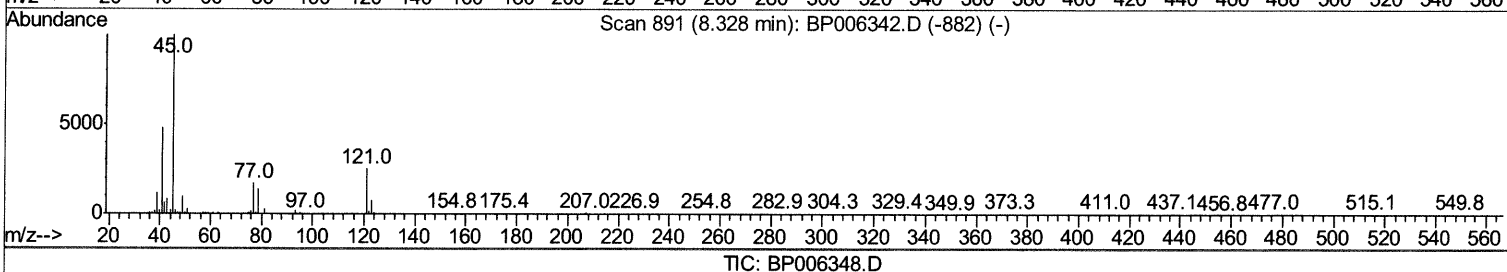
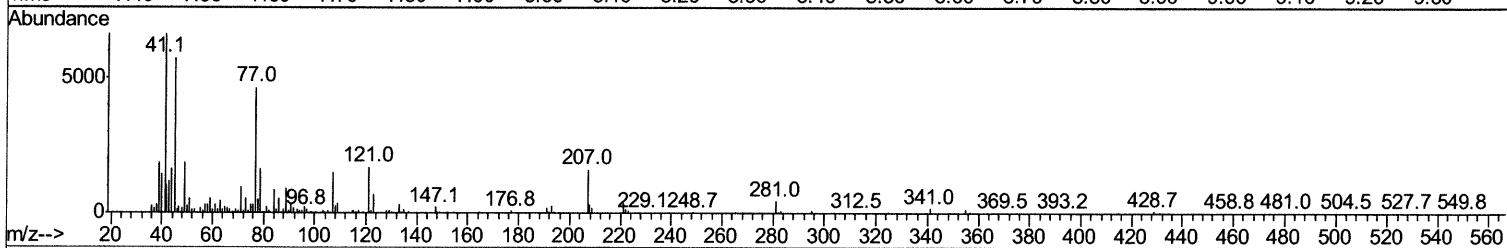
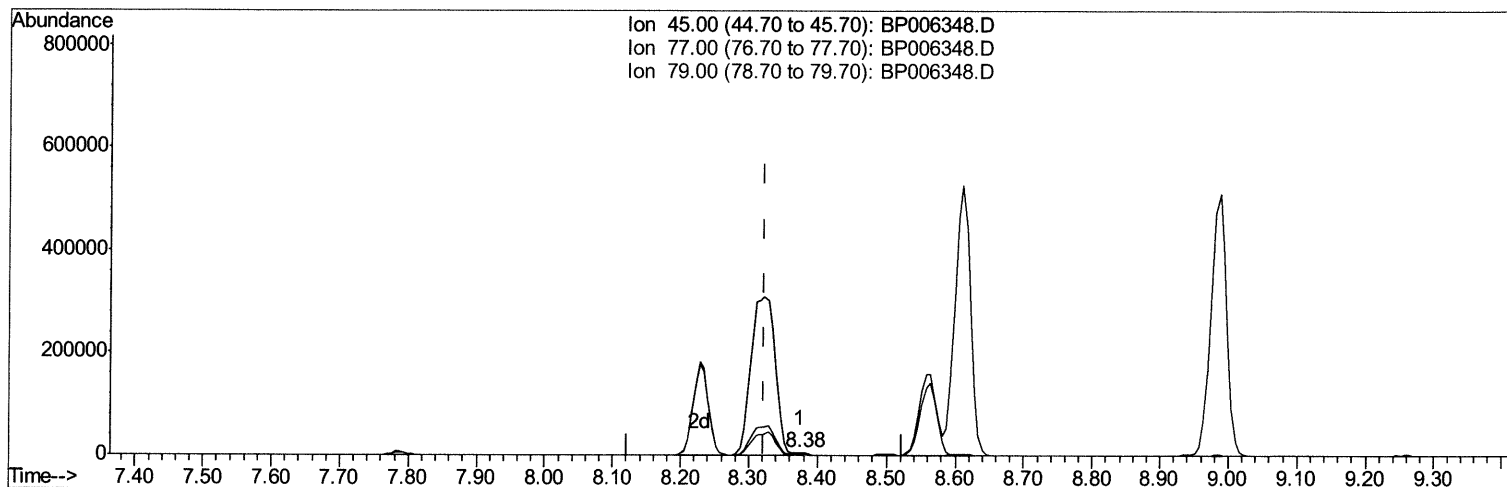
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP072621\
Data File : BP006348.D
Acq On : 26 Jul 2021 20:51
Operator : CG/JU
Sample : PB137859BS
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SLCS859

Manual Integrations
APPROVED

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Quant Time: Jul 27 01:09:51 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jul 24 04:29:12 2021
Response via : Initial Calibration



(14) 2,2'-oxybis(1-Chloropropane)

8.375min (+0.053) 0.33ng/ul

response 6763

Ion	Exp%	Act%
45.00	100	100
77.00	18.20	80.80#
79.00	13.30	28.38#
0.00	0.00	0.00

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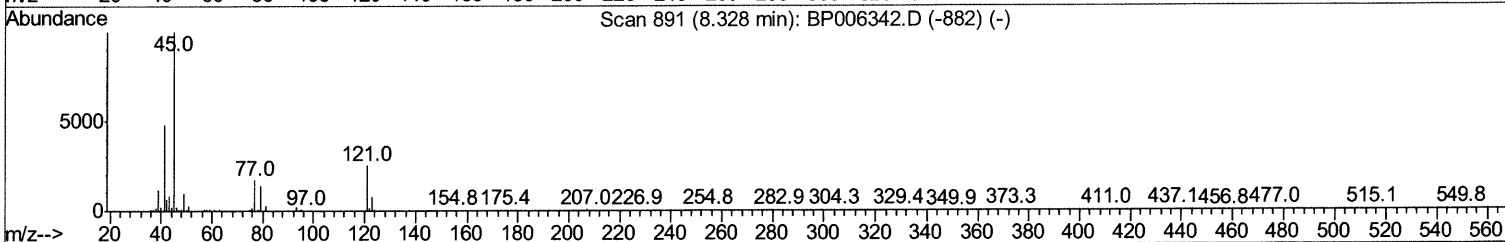
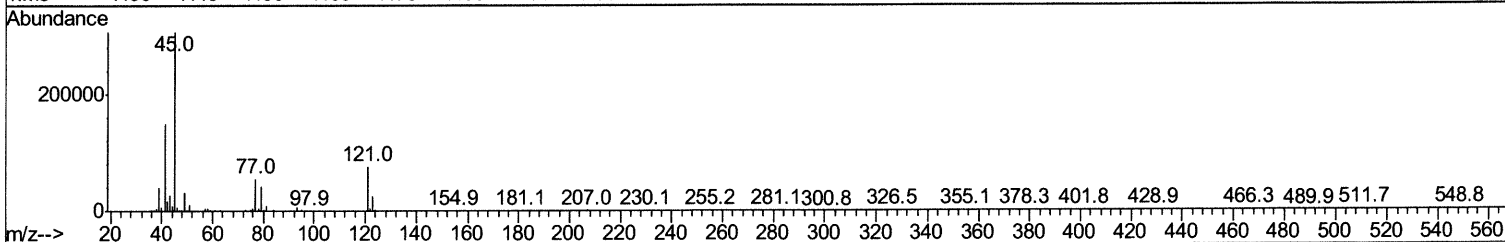
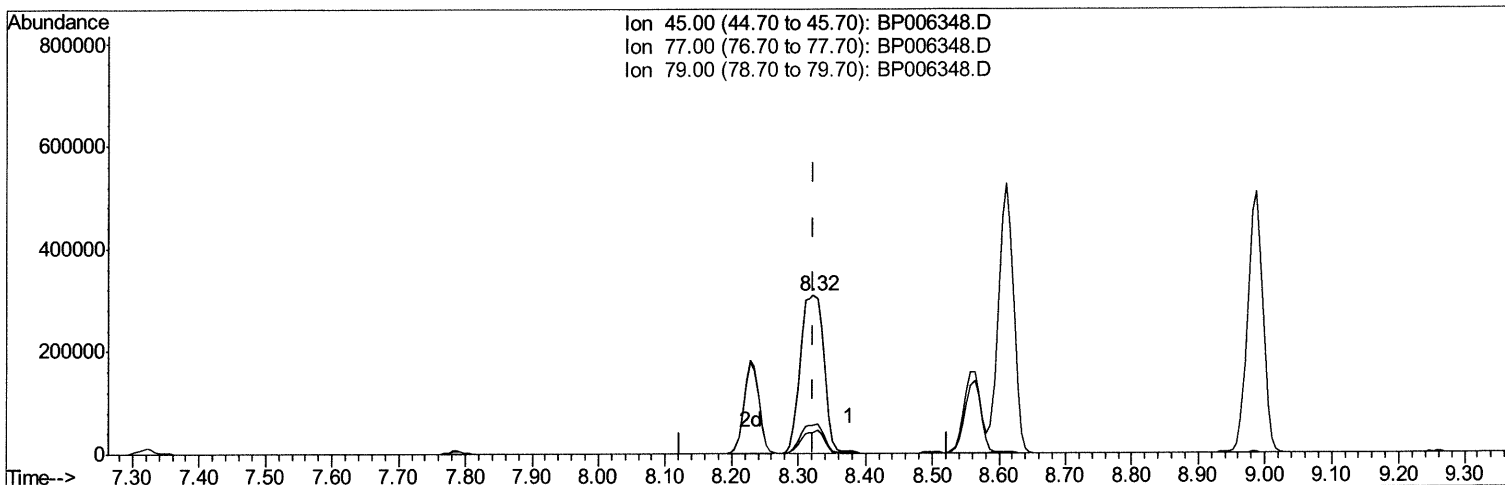
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TIC: BP006348.D

(14) 2,2'-oxybis(1-Chloropropane)

8.322min (-0.000) 37.06ng/ul m 07/29/21 JU

response 758440

Ion	Exp%	Act%
45.00	100	100
77.00	18.20	17.93
79.00	13.30	13.48
0.00	0.00	0.00

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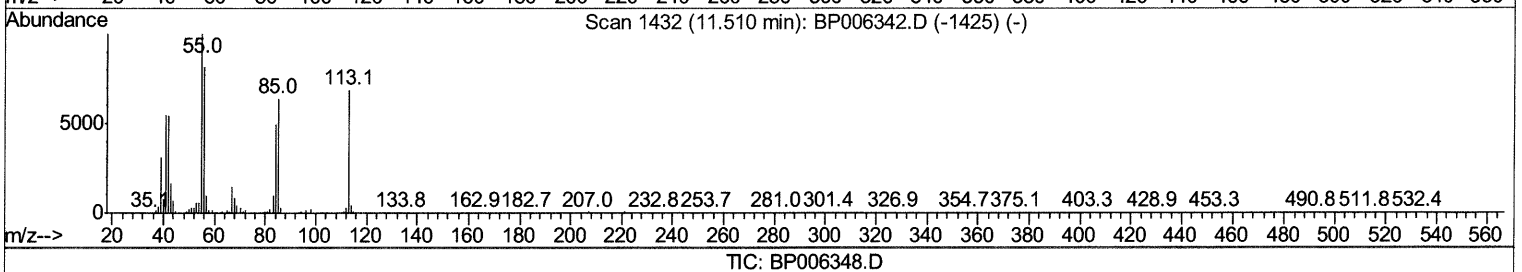
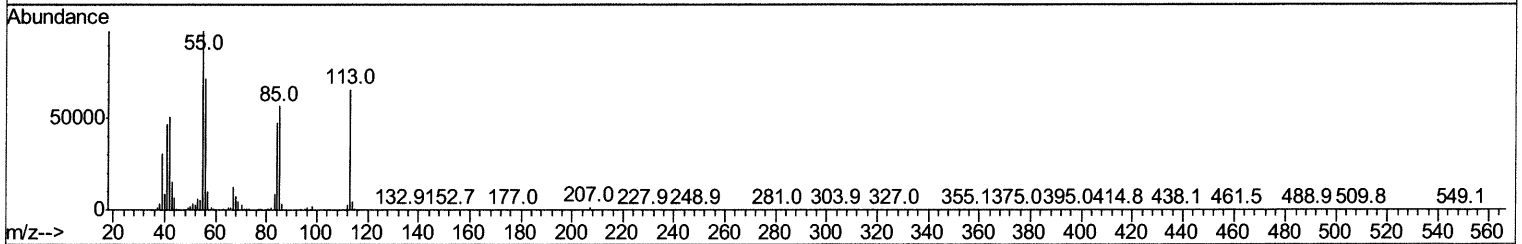
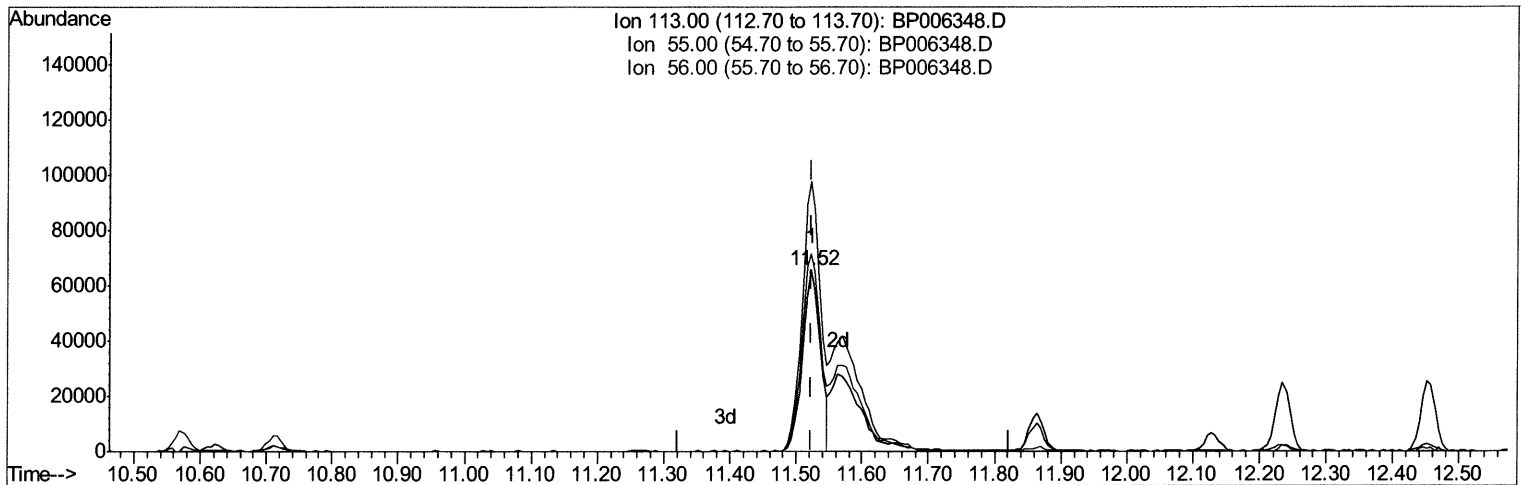
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(34) Caprolactam

11.522min (-0.000) 24.02ng/ul

response 122811

Ion	Exp%	Act%
113.00	100	100
55.00	140.20	148.68
56.00	111.50	108.79
0.00	0.00	0.00

Quantitation Report (Qedit)

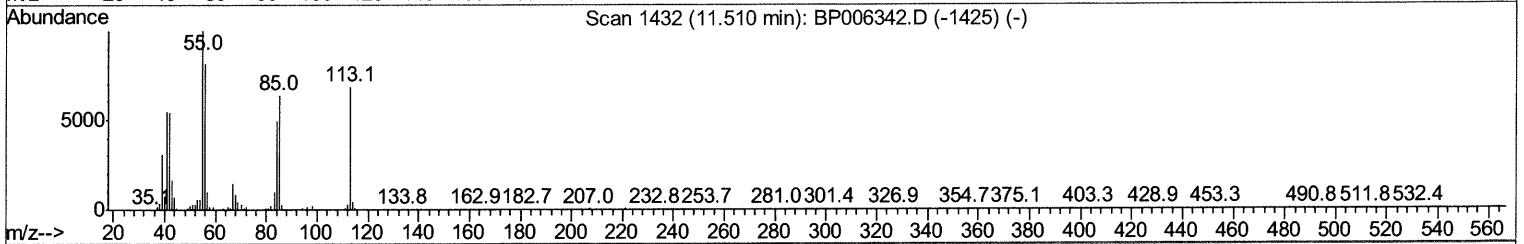
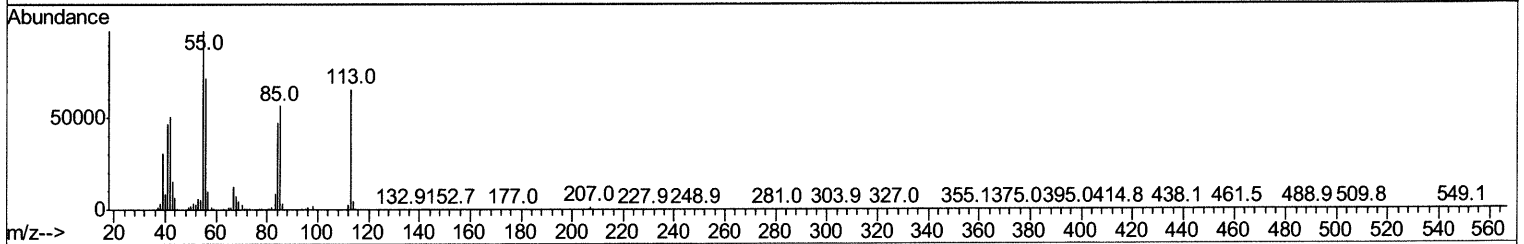
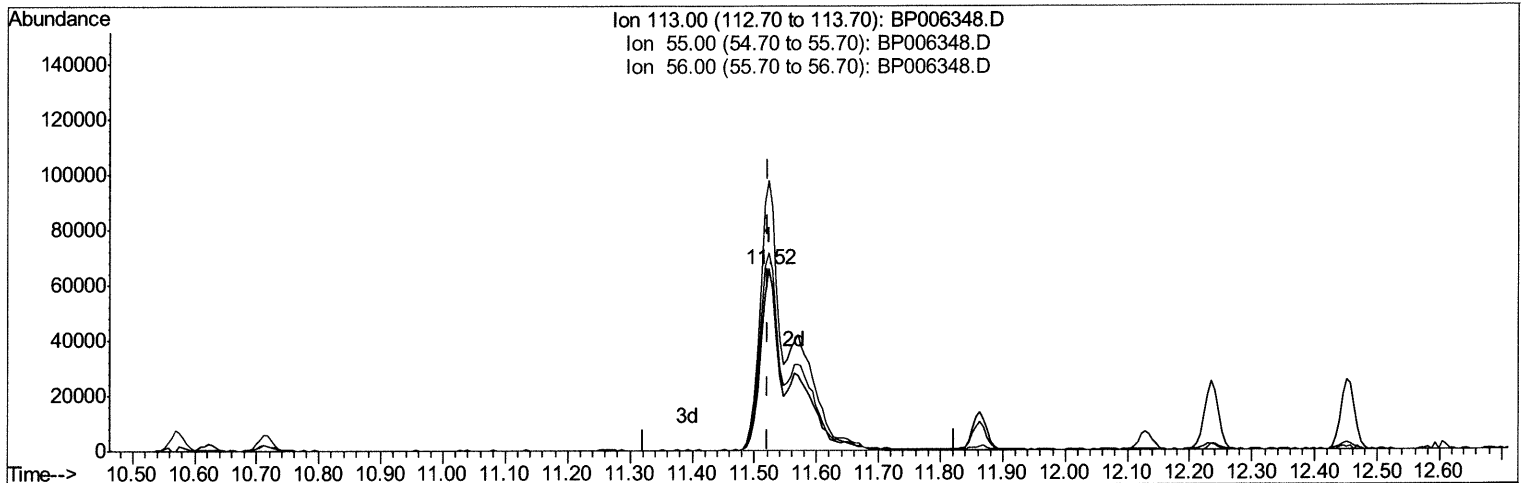
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TIC: BP006348.D

(34) Caprolactam

11.522min (-0.000) 41.85ng/ul m 07/29/2021

response 213992

Ion	Exp%	Act%
113.00	100	100
55.00	140.20	148.68
56.00	111.50	108.79
0.00	0.00	0.00

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 Client Sampled :
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Quant Time: Jul 27 01:36:33 2021
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 Quant Title : SVOA CALIBRATION
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 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	242591	20.00	ng/ul	0.00
20) Naphthalene-d8	10.57	136	1083338	20.00	ng/ul	-0.01
38) Acenaphthene-d10	14.42	164	640758	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.17	188	1248113	20.00	ng/ul	0.00
79) Chrysene-d12	21.26	240	1191128	20.00	ng/ul	0.00
88) Perylene-d12	23.60	264	1209261	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.30	96	42782	6.58	ng/uL	0.00
4) Pyridine-d5	3.70	84	649608	33.67	ng/ul	0.00
7) Phenol-d5	6.96	99	838109	36.79	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.13	67	519835	36.67	ng/ul	0.00
11) 2-Chlorophenol-d4	7.32	132	633532	37.43	ng/ul	0.00
15) 4-Methylphenol-d8	8.50	113	665608	37.15	ng/ul	0.00
21) Nitrobenzene-d5	8.95	128	312795	38.78	ng/ul	0.00
24) 2-Nitrophenol-d4	9.66	143	313664	40.72	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.19	165	568584	37.67	ng/ul	0.00
31) 4-Chloroaniline-d4	10.72	131	870422	35.03	ng/ul	0.00
46) Dimethylphthalate-d6	13.83	166	1780397	39.04	ng/ul	0.00
49) Acenaphthylene-d8	14.10	160	2201686	39.30	ng/ul	0.00
54) 4-Nitrophenol-d4	14.62	143	367901	40.63	ng/ul	0.00
60) Fluorene-d10	15.41	176	1468902	38.31	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.53	200	269601	38.75	ng/ul	0.00
73) Anthracene-d10	17.26	188	2250095	39.94	ng/ul	0.00
81) Pyrene-d10	19.50	212	2505102	40.89	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.46	264	2435673	39.64	ng/ul	0.00

Target Compounds

					Ovalue	
2) 1,4-Dioxane	3.33	88	88852	12.949	ng/uL	96
5) Pyridine	3.72	79	661827	32.893	ng/ul	98
6) Benzaldehyde	6.93	77	508573	40.912	ng/ul	97
8) Phenol	6.99	94	852850	36.620	ng/ul	97
10) Bis(2-Chloroethyl)ether	7.22	93	651728	35.576	ng/ul	98
12) 2-Chlorophenol	7.35	128	640309	37.312	ng/ul	98
13) 2-Methylphenol	8.23	108	615947	36.243	ng/ul	99
14) 2,2'-oxybis(1-Chloropropan	8.32	45	758440m	37.059	ng/ul	>
16) Acetophenone	8.61	105	1007530	36.568	ng/ul	97
17) N-Nitroso-di-n-propylamine	8.60	70	564083	39.134	ng/ul	97
18) 4-Methylphenol	8.56	108	678720	37.192	ng/ul	99
19) Hexachloroethane	8.86	117	259898	37.024	ng/ul	99
22) Nitrobenzene	8.99	77	813251	38.858	ng/ul	99
23) Isophorone	9.52	82	1489900	39.821	ng/ul	99
25) 2-Nitrophenol	9.69	139	336573	39.627	ng/ul	97
26) 2,4-Dimethylphenol	9.76	107	720884	35.789	ng/ul	98
27) Bis(2-Chloroethoxy)methane	10.00	93	901104	37.504	ng/ul	100
29) 2,4-Dichlorophenol	10.22	162	559160	37.761	ng/ul	99
30) Naphthalene	10.62	128	2023473	36.000	ng/ul	100
32) 4-Chloroaniline	10.74	127	835845	34.147	ng/ul	99
33) Hexachlorobutadiene	10.91	225	319723	35.392	ng/ul	99
34) Caprolactam	11.52	113	213992m	41.852	ng/ul	>

07/29/21 JU

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
35) 4-Chloro-3-methylphenol	11.86	107	715695	40.487	ng/ul	98
36) 2-Methylnaphthalene	12.23	142	1425470	37.012	ng/ul	100
37) 1-Methylnaphthalene	12.45	142	1395570	35.999	ng/ul	99
39) 1,2,4,5-Tetrachlorobenzene	12.60	216	621654	36.001	ng/ul	99
40) Hexachlorocyclopentadiene	12.58	237	347920	30.973	ng/ul	99
41) 2,4,6-Trichlorophenol	12.85	196	421857	39.426	ng/ul	98
42) 2,4,5-Trichlorophenol	12.92	196	462510	39.507	ng/ul	99
43) 1,1'-Biphenyl	13.25	154	1771327	36.878	ng/ul	99
44) 2-Chloronaphthalene	13.29	162	1325432	36.548	ng/ul	98
45) 2-Nitroaniline	13.50	65	475872	44.421	ng/ul	98
47) Dimethylphthalate	13.89	163	1749013	39.072	ng/ul	99
48) 2,6-Dinitrotoluene	14.00	165	350398	42.817	ng/ul	95
50) Acenaphthylene	14.13	152	2074407	38.019	ng/ul	100
51) 3-Nitroaniline	14.32	138	377267	39.345	ng/ul	96
52) Acenaphthene	14.48	153	1505660	37.611	ng/ul	99
53) 2,4-Dinitrophenol	14.53	184	183547	37.817	ng/ul	97
55) 4-Nitrophenol	14.63	109	318002	43.418	ng/ul	95
56) Dibenzofuran	14.82	168	2010098	37.441	ng/ul	97
57) 2,4-Dinitrotoluene	14.78	165	508856	43.512	ng/ul	93
58) 2,3,4,6-Tetrachlorophenol	15.04	232	372215	40.770	ng/ul	97
59) Diethylphthalate	15.26	149	1881582	40.582	ng/ul	100
61) Fluorene	15.46	166	1682500	38.099	ng/ul	100
62) 4-Chlorophenyl-phenylether	15.47	204	766520	37.231	ng/ul	99
63) 4-Nitroaniline	15.49	138	411974	43.745	ng/ul	97
66) 4,6-Dinitro-2-methylphenol	15.55	198	267170	39.279	ng/ul#	96
67) N-Nitrosodiphenylamine	15.68	169	1444931	38.970	ng/ul	100
68) 4-Bromophenyl-phenylether	16.36	248	449824	45.003	ng/ul	97
69) Hexachlorobenzene	16.46	284	519515	38.755	ng/ul	97
70) Atrazine	16.65	200	516757	40.273	ng/ul	99
71) Pentachlorophenol	16.81	266	329339	39.081	ng/ul	99
72) Phenanthrene	17.21	178	2626052	39.183	ng/ul	99
74) Anthracene	17.30	178	2647806	38.964	ng/ul	99
75) 1,2,3,4-Tetrachlorobenzene	13.21	216	625874	36.583	ng/uL	99
76) Pentachlorobenzene	14.73	250	621019	37.235	ng/uL	98
77) Carbazole	17.57	167	2483136	39.945	ng/ul	99
78) Di-n-butylphthalate	18.14	149	3231563	43.496	ng/ul	99
80) Fluoranthene	19.18	202	2977988	39.448	ng/ul	99
82) Pyrene	19.53	202	3051879	38.959	ng/ul	98
83) Butylbenzylphthalate	20.43	149	1437118	44.712	ng/ul	100
84) 3,3'-Dichlorobenzidine	21.19	252	997385	38.003	ng/ul	99
85) Benzo(a)anthracene	21.25	228	2905678	39.368	ng/ul	100
86) Bis(2-ethylhexyl)phthalate	21.19	149	2241334	44.538	ng/ul	99
87) Chrysene	21.30	228	2803527	39.627	ng/ul	99
89) Di-n-octyl phthalate	22.10	149	3733392	44.302	ng/ul	100
90) Benzo(b)fluoranthene	22.90	252	3106829	40.449	ng/ul	99
91) Benzo(k)fluoranthene	22.94	252	2855105	39.165	ng/ul	99
93) Benzo(a)pyrene	23.50	252	2694070	40.127	ng/ul	100
94) Indeno(1,2,3-cd)pyrene	26.04	276	3400667	39.570	ng/ul	97
95) Dibenzo(a,h)anthracene	26.06	278	2879456	39.559	ng/ul	99
96) Benzo(g,h,i)perylene	26.78	276	2819393	39.706	ng/ul	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
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(#) = qualifier out of range (m) = manual integration (+) = signals summed