

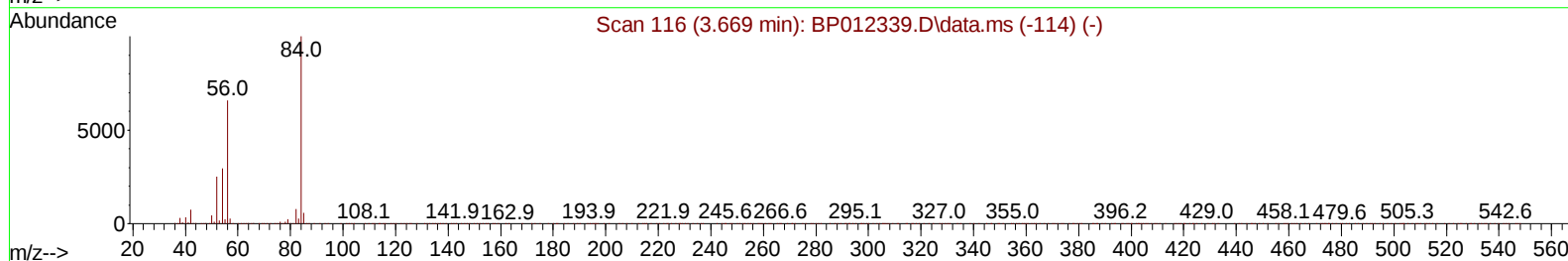
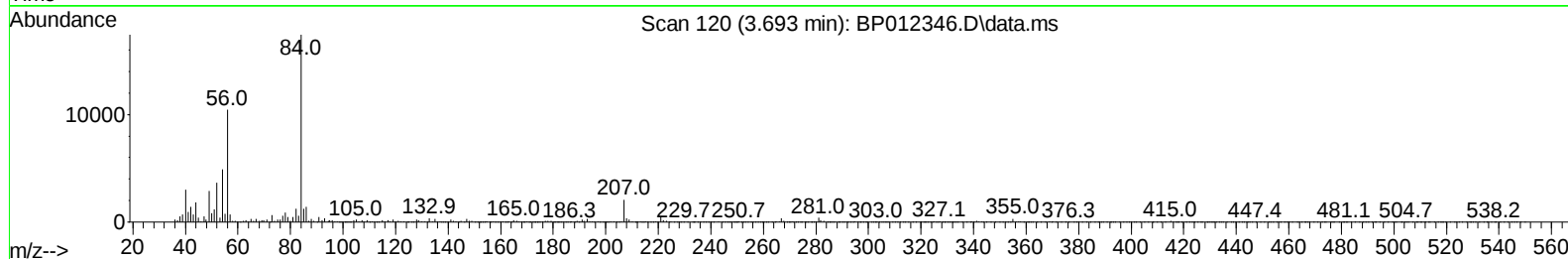
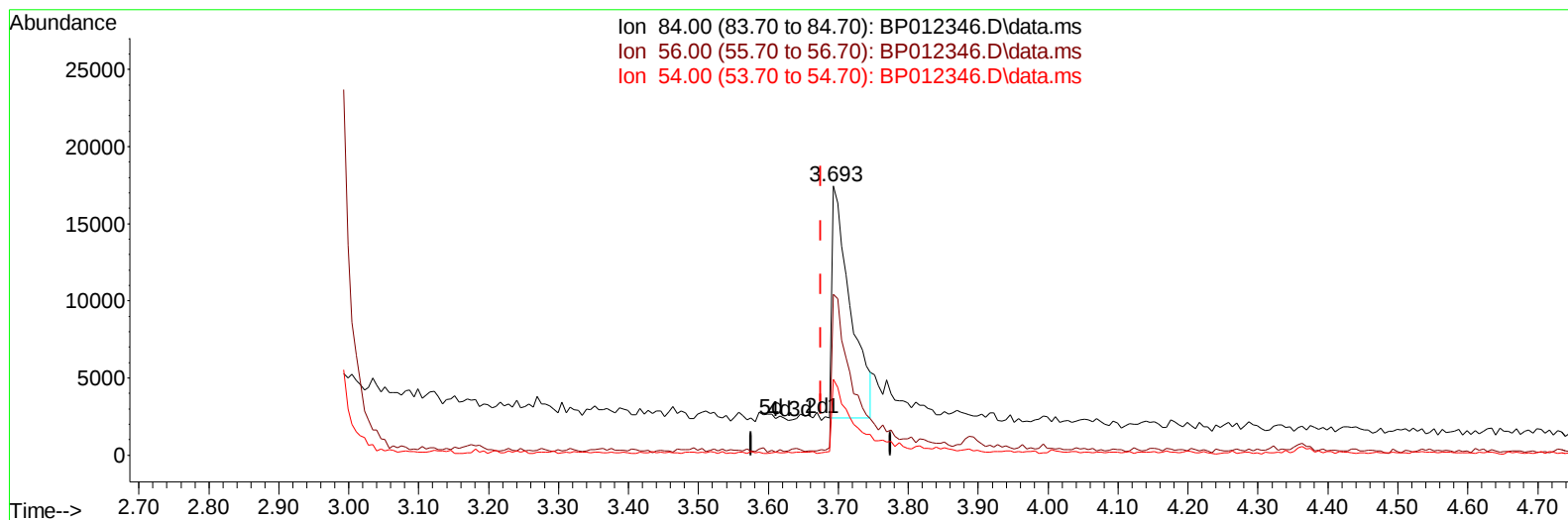
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012346.D
Acq On : 27 Oct 2022 16:39
Operator : CG/JU
Sample : N5015-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C1BG4

Manual IntegrationsAPPROVED

Quant Time: Oct 27 23:22:06 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 26 23:40:20 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/28/2022
Supervised By :mohammad ahmed 10/31/2022



TIC: BP012346.D\data.ms

(4) Pyridine-d5 (S)

3.693min (+ 0.018) 1.33 ng/u1

response 27420

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	59.87
54.00	29.80	28.20
0.00	0.00	0.00

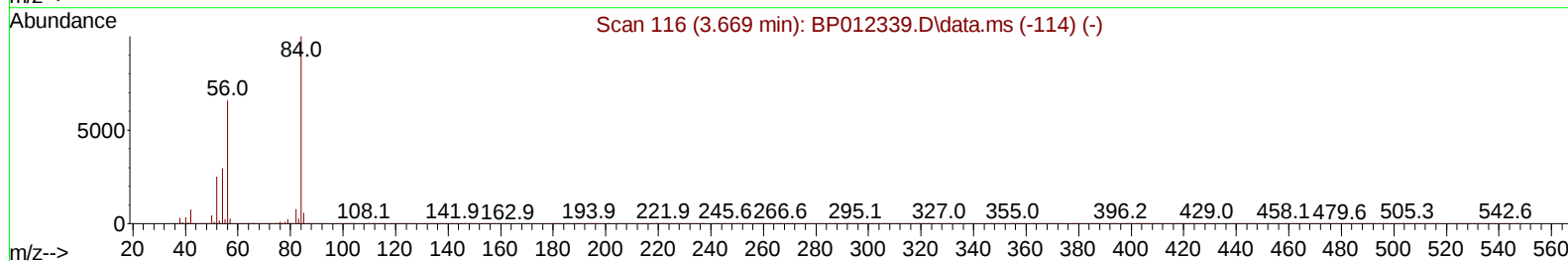
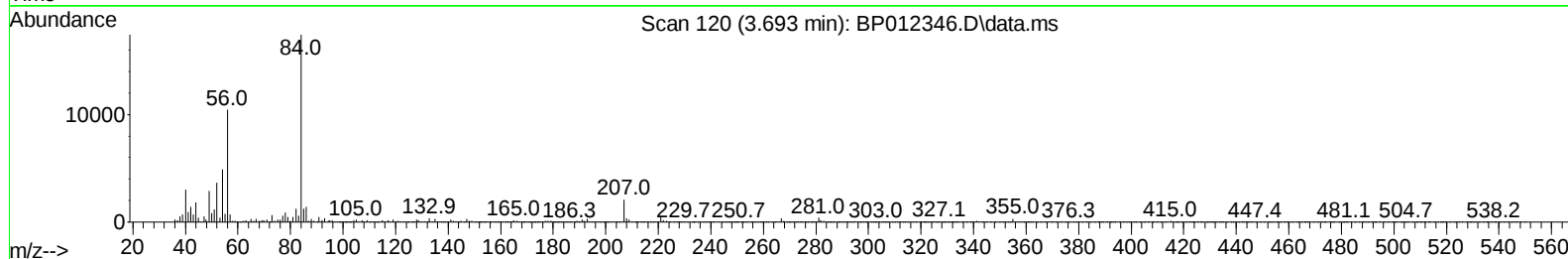
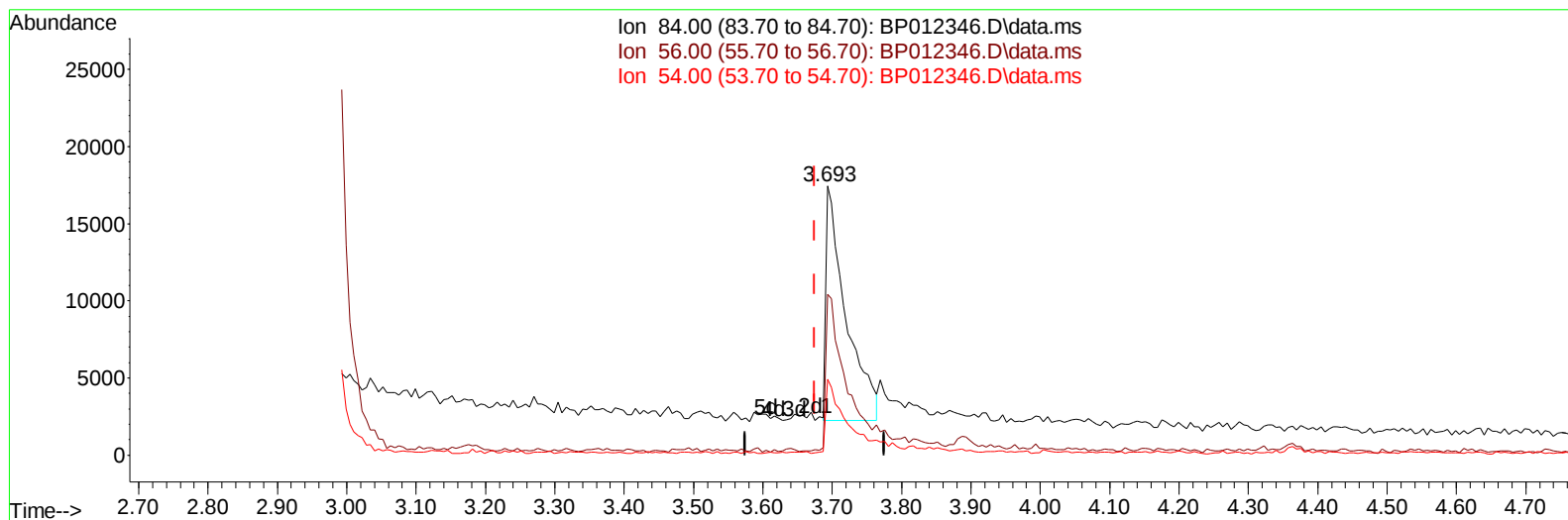
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TIC: BP012346.D\data.ms

(4) Pyridine-d5 (S)

3.693min (+ 0.018) 1.48 ng/u1 m

response 30497

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	59.87
54.00	29.80	28.20
0.00	0.00	0.00

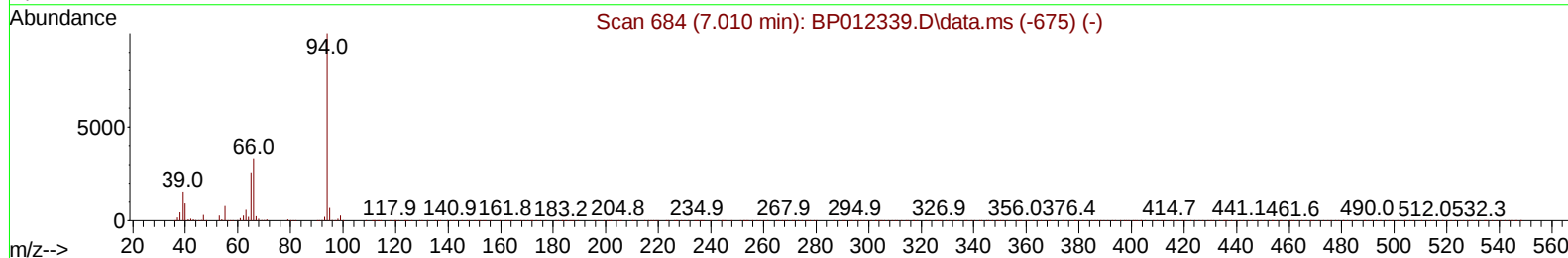
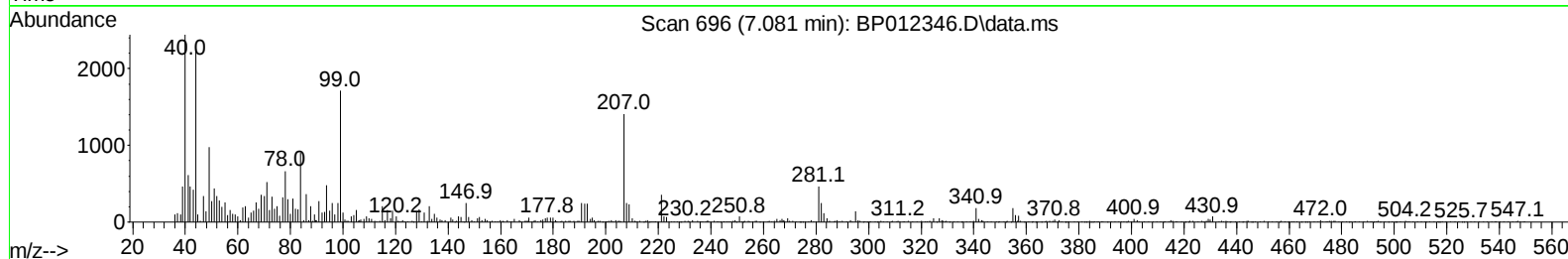
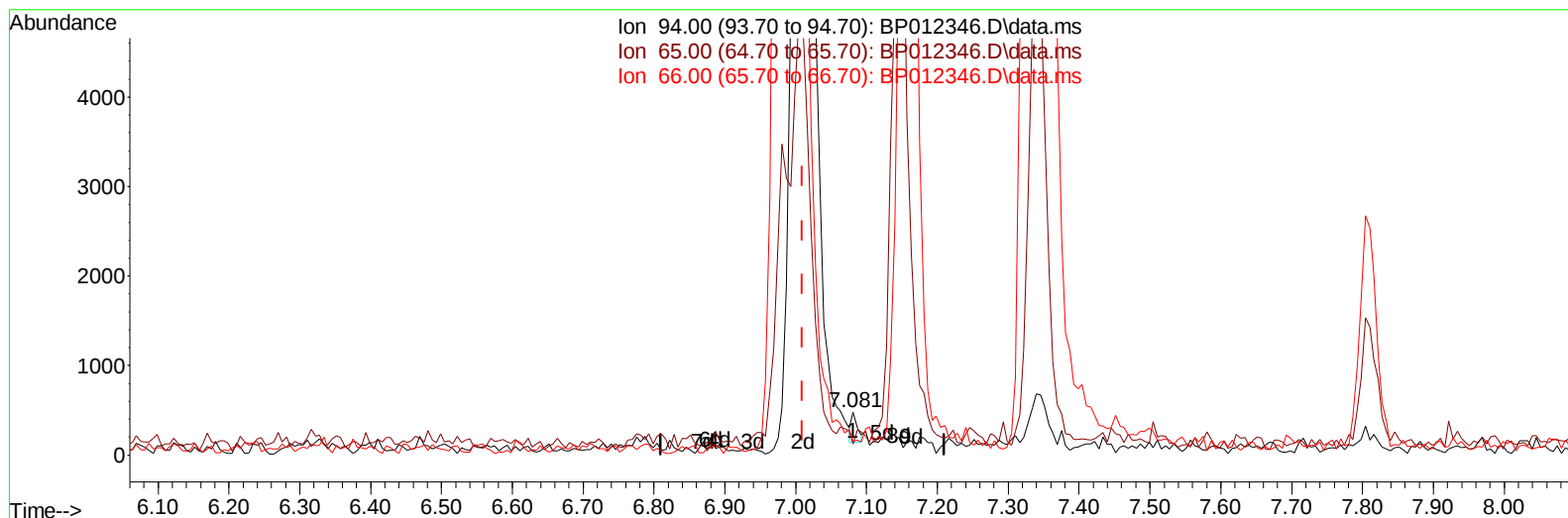
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TIC: BP012346.D\data.ms

(8) Phenol

7.081min (+ 0.071) 0.01 ng/u1

response 202

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	26.00	26.04
66.00	32.70	31.25
0.00	0.00	0.00

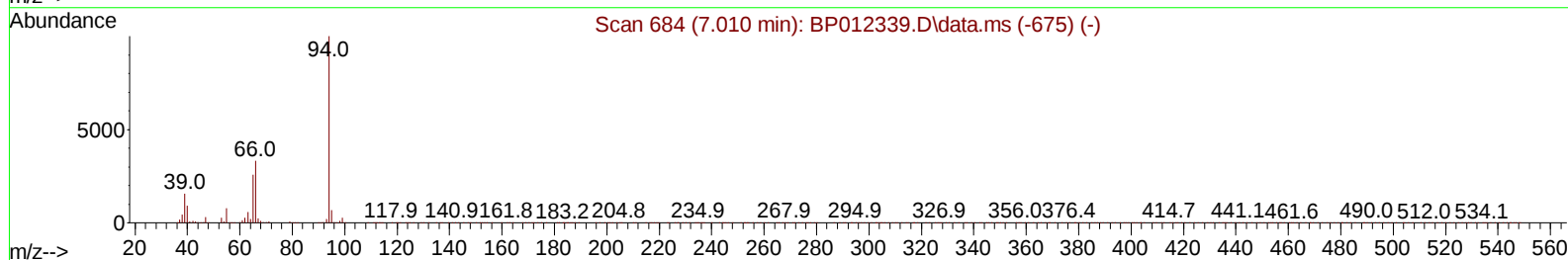
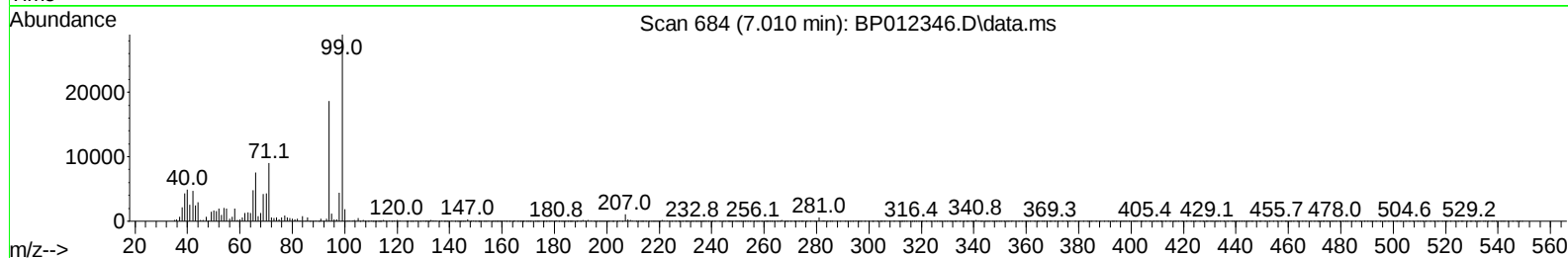
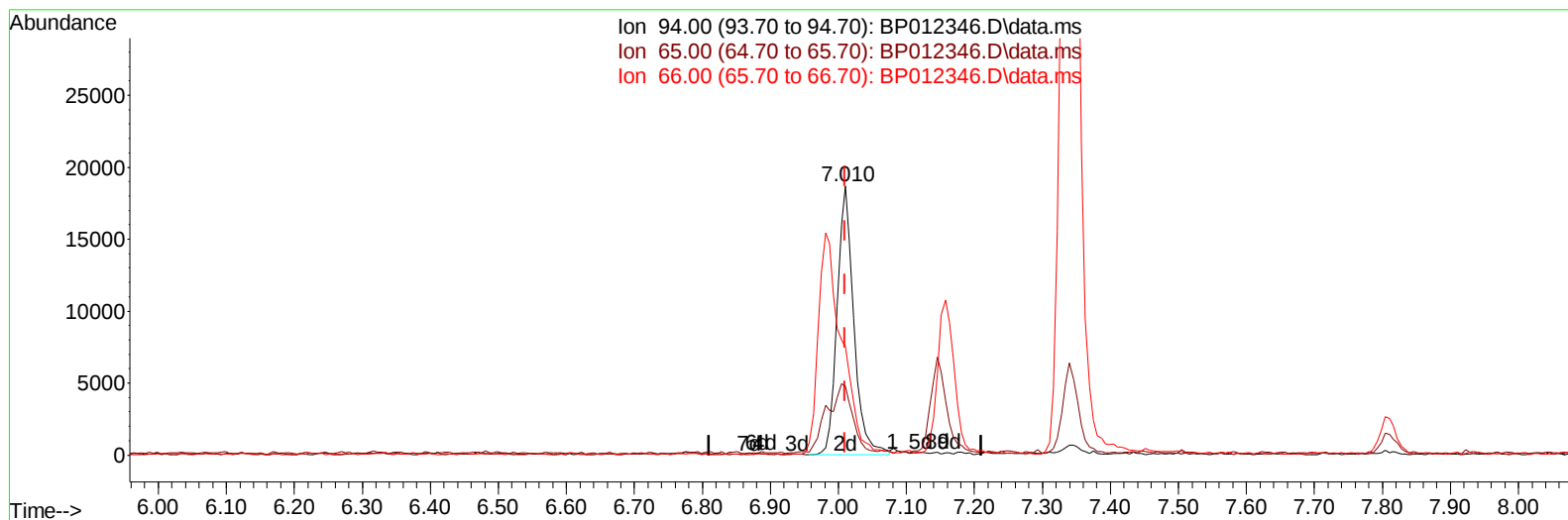
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TIC: BP012346.D\data.ms

(8) Phenol

7.010min (+ 0.000) 1.18 ng/ul m

response 31788

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	26.00	25.65
66.00	32.70	40.41#
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	288965	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	1307970	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	887161	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.222	188	1933091	20.000	ng/ul	0.00
79) Chrysene-d12	21.315	240	2034093	20.000	ng/ul	0.00
88) Perylene-d12	23.715	264	1911211	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	31437	4.380	ng/uL	0.00
4) Pyridine-d5	3.693	84	30497m	1.479	ng/ul	0.02
7) Phenol-d5	6.981	99	612437	23.630	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.146	67	389301	25.161	ng/ul	0.00
11) 2-Chlorophenol-d4	7.340	132	493238	26.060	ng/ul	0.00
15) 4-Methylphenol-d8	8.528	113	429325	20.944	ng/ul	0.00
21) Nitrobenzene-d5	8.981	128	256240	30.210	ng/ul	0.00
24) 2-Nitrophenol-d4	9.698	143	276802	31.442	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.240	165	491226	25.747	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	650722	22.906	ng/ul	0.00
46) Dimethylphthalate-d6	13.875	166	1622748	26.720	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	1862466	26.836	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	190669	16.589	ng/ul	0.00
60) Fluorene-d10	15.457	176	1452053	27.926	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	163951	16.290	ng/ul	0.00
73) Anthracene-d10	17.322	188	2175408	26.118	ng/ul	0.00
81) Pyrene-d10	19.563	212	2780488	27.226	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.557	264	2531978	27.113	ng/ul	0.00
Target Compounds						
8) Phenol	7.010	94	31788m	1.180	ng/ul	
80) Fluoranthene	19.233	202	168147	1.334	ng/ul	99
82) Pyrene	19.592	202	144248	1.104	ng/ul	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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