

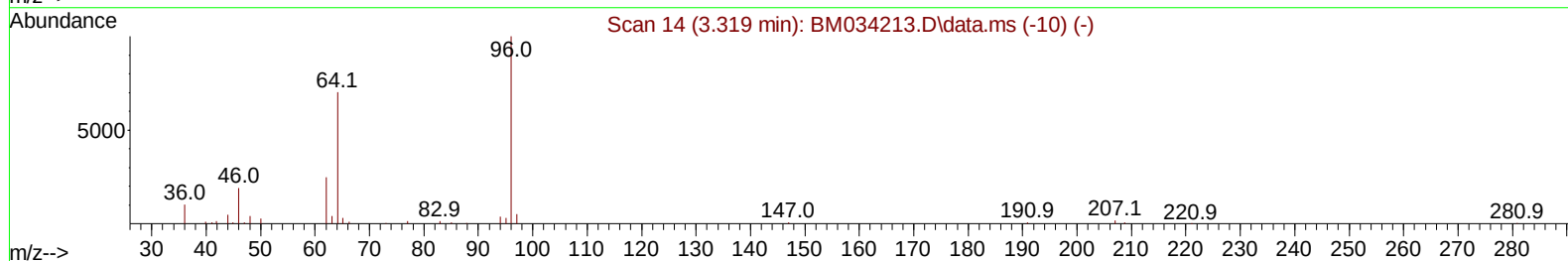
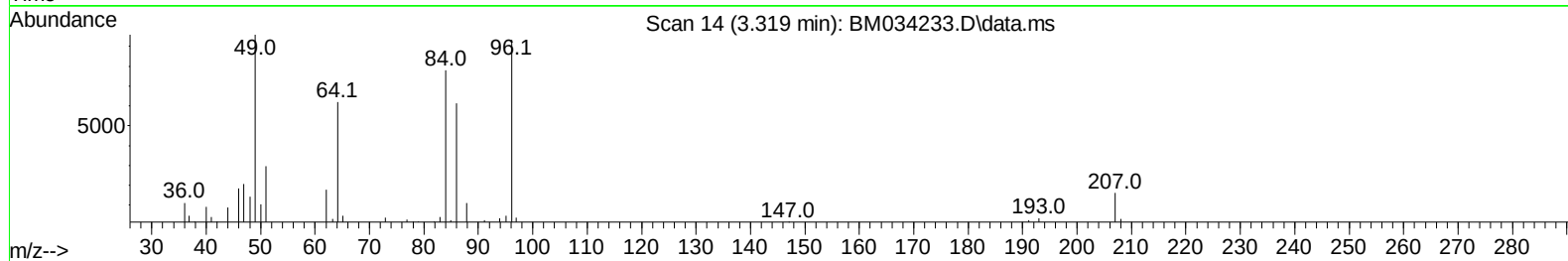
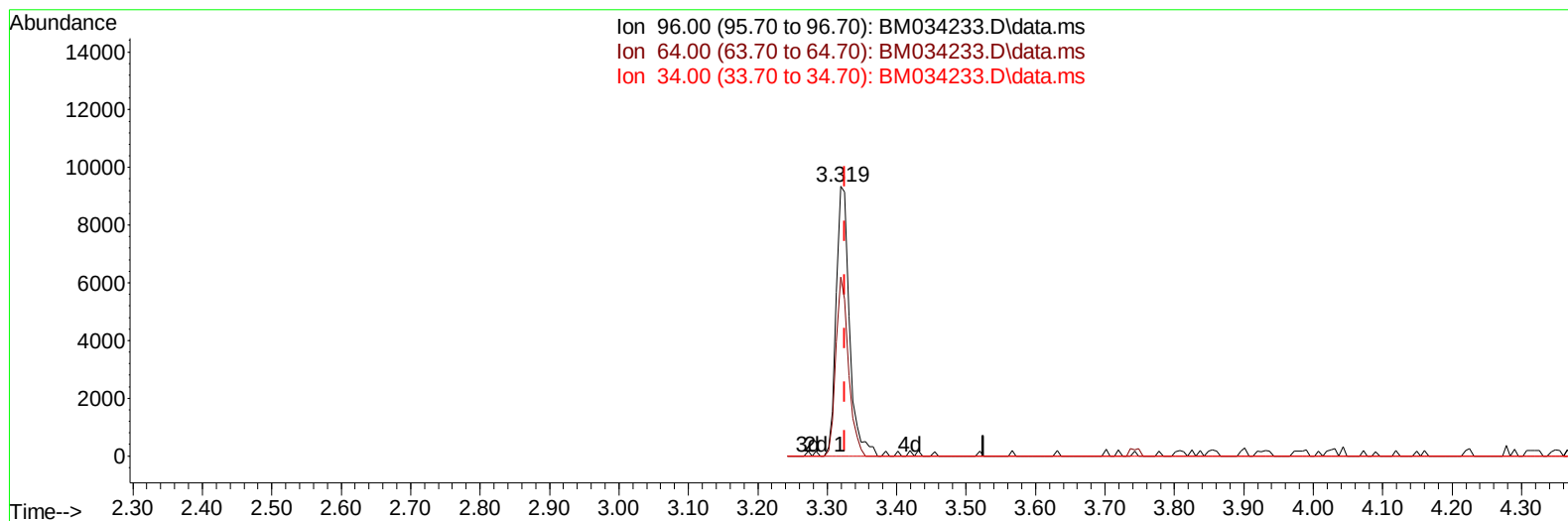
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031122\
 Data File : BM034233.D
 Acq On : 11 Mar 2022 22:13
 Operator : CG/JU
 Sample : N1816-15
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBSE0

Manual IntegrationsAPPROVED

Quant Time: Mar 11 23:42:57 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM031022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 11 02:00:05 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/14/2022
 Supervised By :mohammad ahmed 03/14/2022



TIC: BM034233.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.319min (-0.006) 3.81 ng/uL

response 12497

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	68.90	66.48
34.00	0.00	0.00
0.00	0.00	0.00

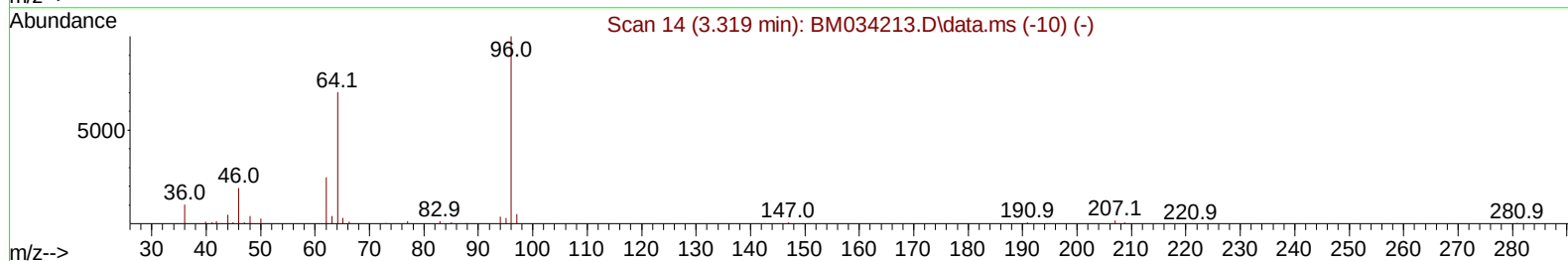
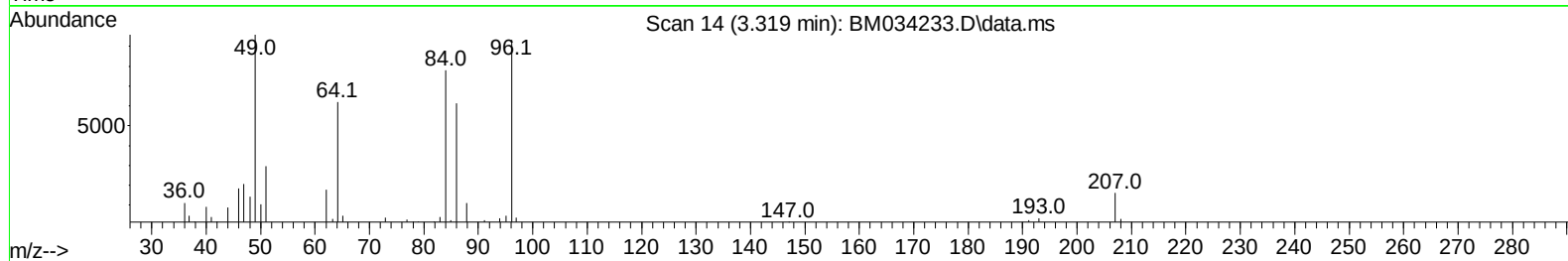
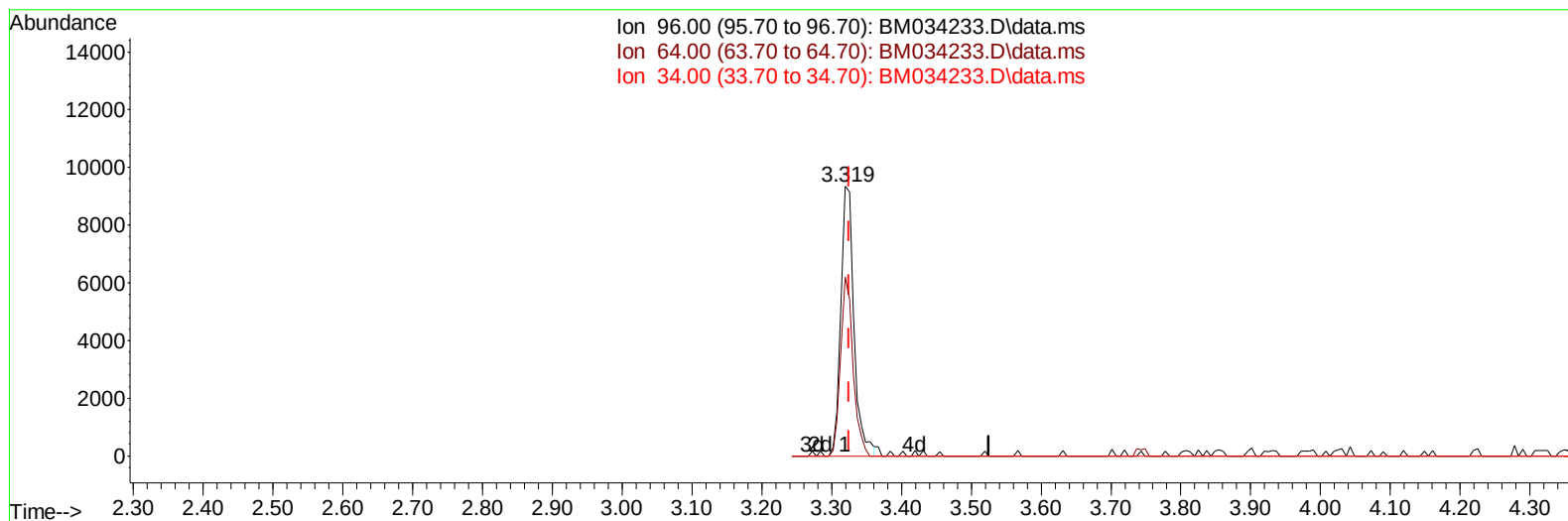
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TIC: BM034233.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.319min (-0.006) 3.78 ng/uL m

response 12380

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	68.90	66.48
34.00	0.00	0.00
0.00	0.00	0.00

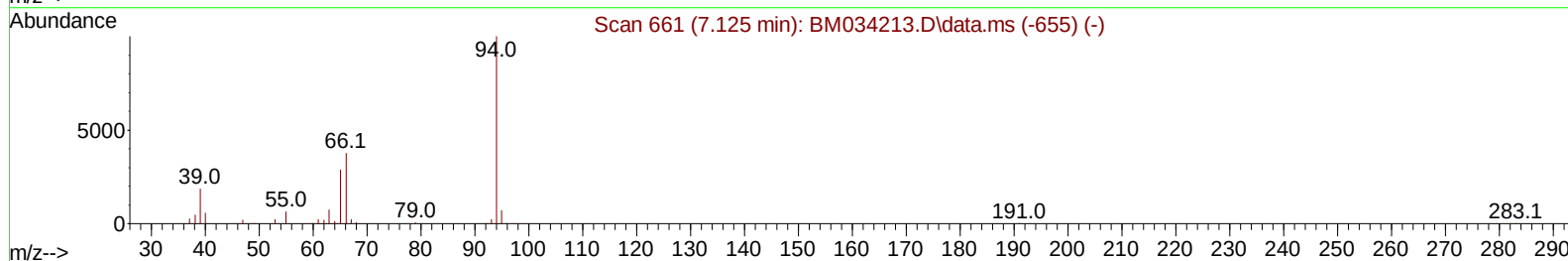
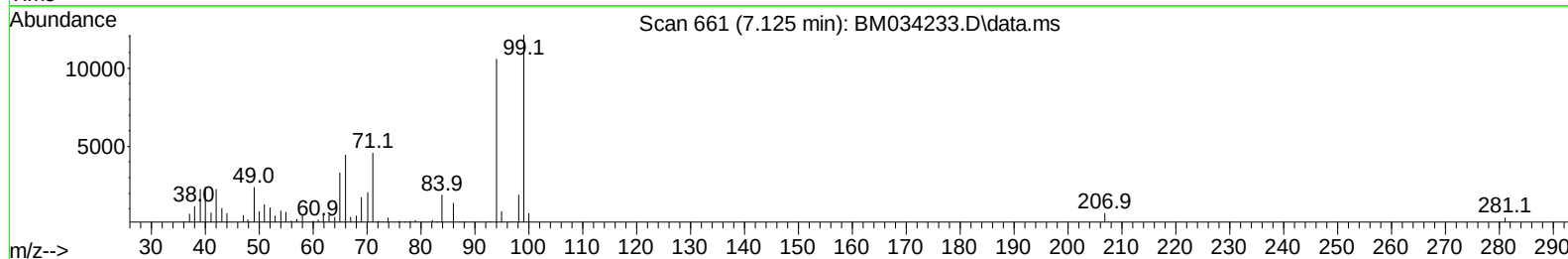
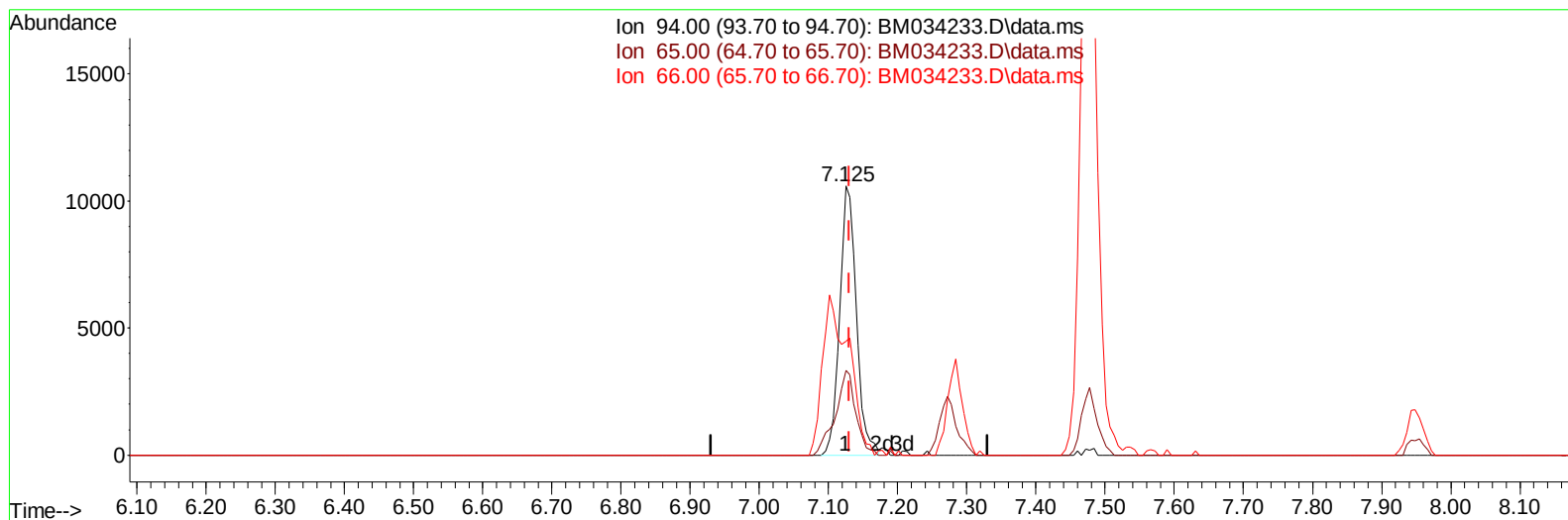
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TIC: BM034233.D\data.ms

(8) Phenol

7.125min (-0.006) 1.88 ng/u1

response 17779

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	28.60	31.42
66.00	42.10	42.24
0.00	0.00	0.00

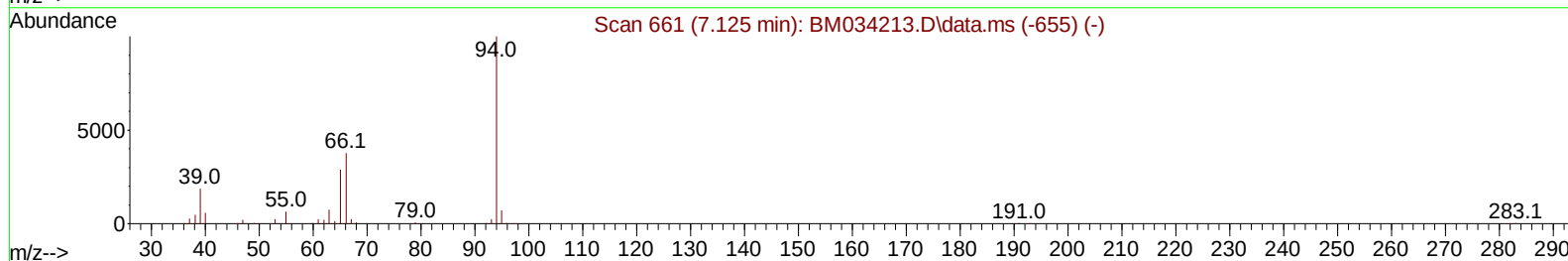
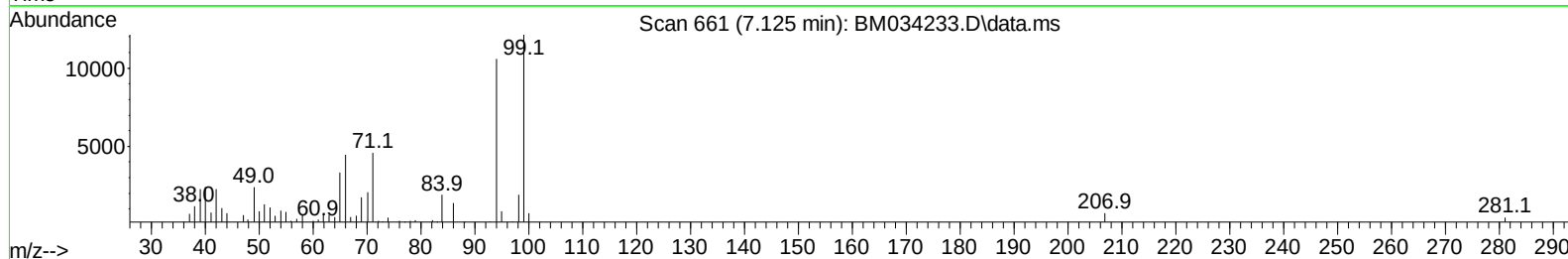
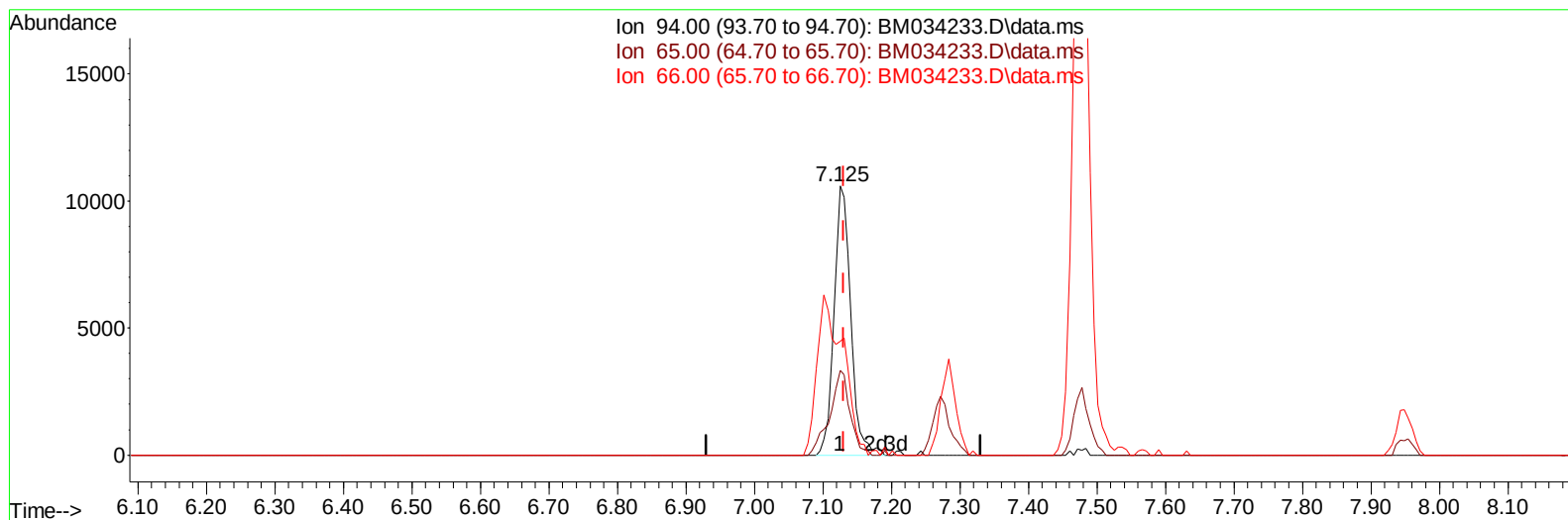
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TIC: BM034233.D\data.ms

(8) Phenol

7.125min (-0.006) 1.90 ng/ul m

response 17955

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	28.60	31.42
66.00	42.10	42.24
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	7.948	152	143133	20.000	ng/ul	0.00
20) Naphthalene-d8	10.766	136	567770	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.589	164	378801	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.330	188	743785	20.000	ng/ul	0.00
79) Chrysene-d12	21.500	240	693920	20.000	ng/ul	0.00
88) Perylene-d12	23.918	264	731711	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	12380m	3.775	ng/uL	0.00
4) Pyridine-d5	3.743	84	124287	15.920	ng/ul	0.00
7) Phenol-d5	7.101	99	176161	18.391	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.272	67	109747	19.530	ng/ul	0.00
11) 2-Chlorophenol-d4	7.478	132	164523	19.875	ng/ul	0.00
15) 4-Methylphenol-d8	8.654	113	152469	18.612	ng/ul	0.00
21) Nitrobenzene-d5	9.107	128	74194	18.218	ng/ul	0.00
24) 2-Nitrophenol-d4	9.837	143	86610	19.179	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.378	165	171462	18.939	ng/ul	0.00
31) 4-Chloroaniline-d4	10.889	131	205259	18.231	ng/ul	0.00
46) Dimethylphthalate-d6	13.995	166	579303	21.018	ng/ul	0.00
49) Acenaphthylene-d8	14.283	160	691534	19.754	ng/ul	0.00
54) 4-Nitrophenol-d4	14.766	143	48450	11.285	ng/ul	0.00
60) Fluorene-d10	15.577	176	501077	20.768	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.689	200	61108	12.985	ng/ul	0.00
73) Anthracene-d10	17.430	188	728713	20.743	ng/ul	0.00
81) Pyrene-d10	19.712	212	870386	23.040	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.759	264	766605	21.236	ng/ul	0.00
Target Compounds						
8) Phenol	7.125	94	17955m	1.900	ng/ul	Qvalue

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

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