

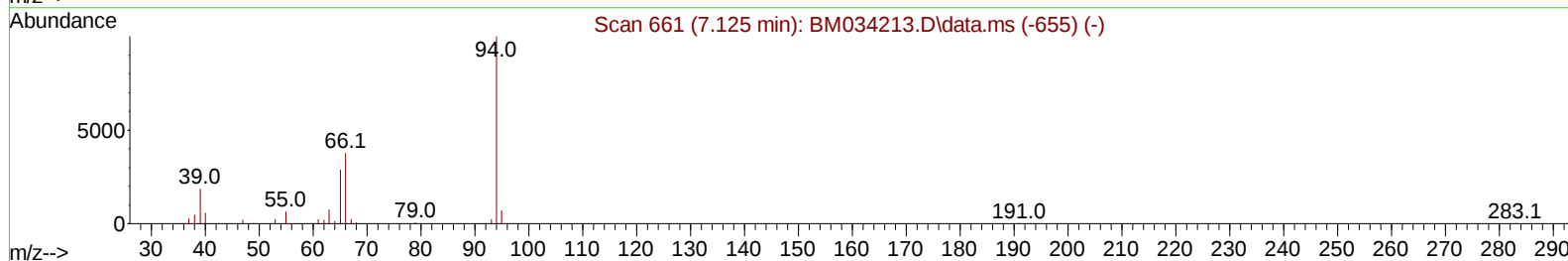
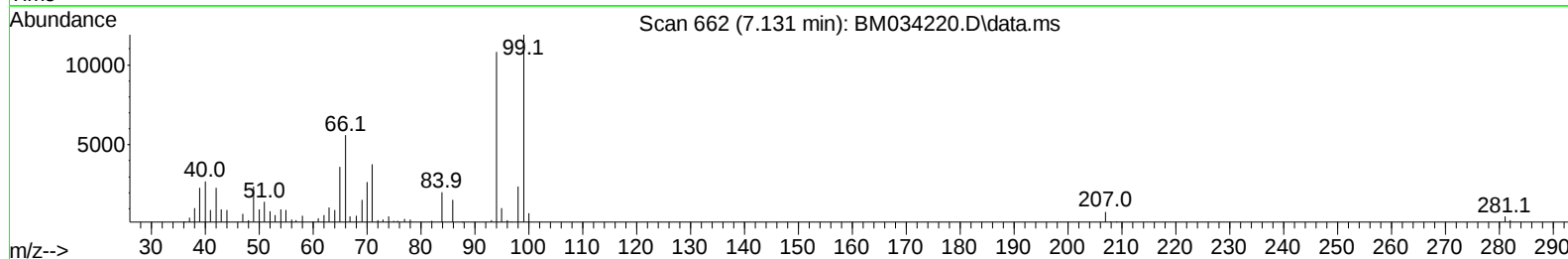
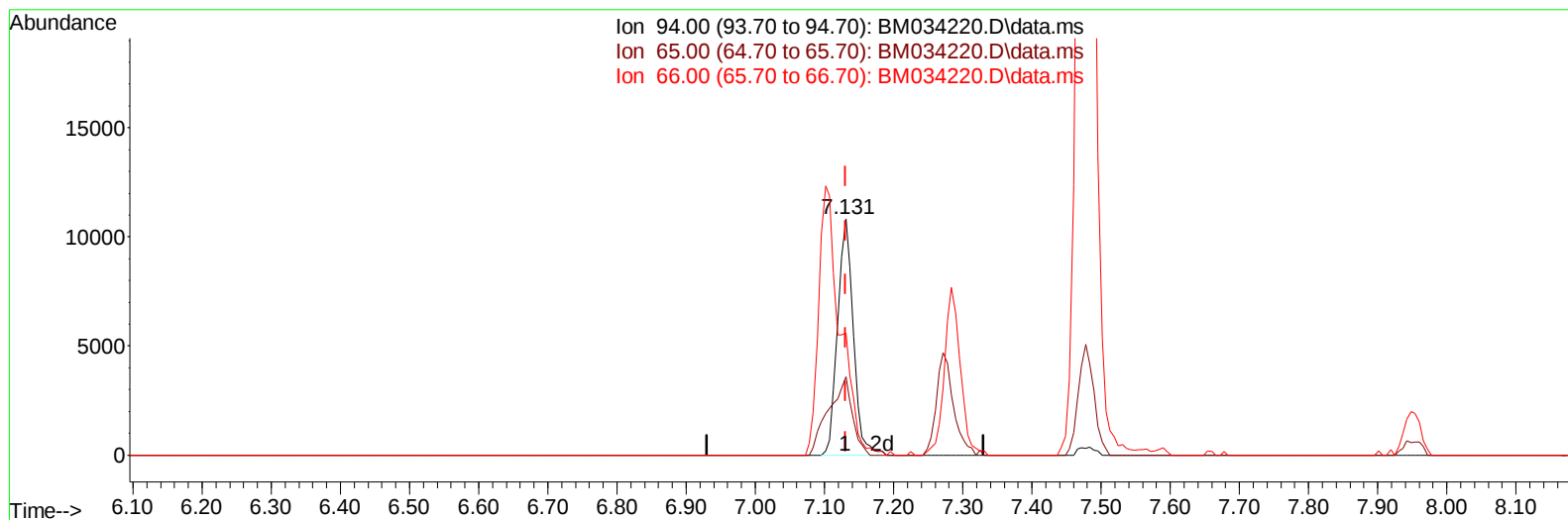
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031122\
 Data File : BM034220.D
 Acq On : 11 Mar 2022 13:46
 Operator : CG/JU
 Sample : N1816-17
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBSE2

Manual IntegrationsAPPROVED

Quant Time: Mar 11 23:40:12 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: BM034220.D\data.ms

(8) Phenol

7.131min (+ 0.000) 1.73 ng/u1

response 17082

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	28.60	33.28
66.00	42.10	51.63#
0.00	0.00	0.00

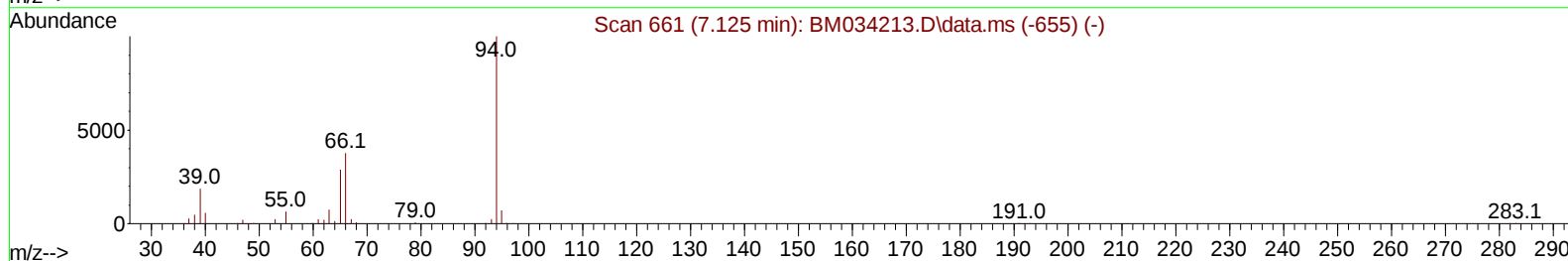
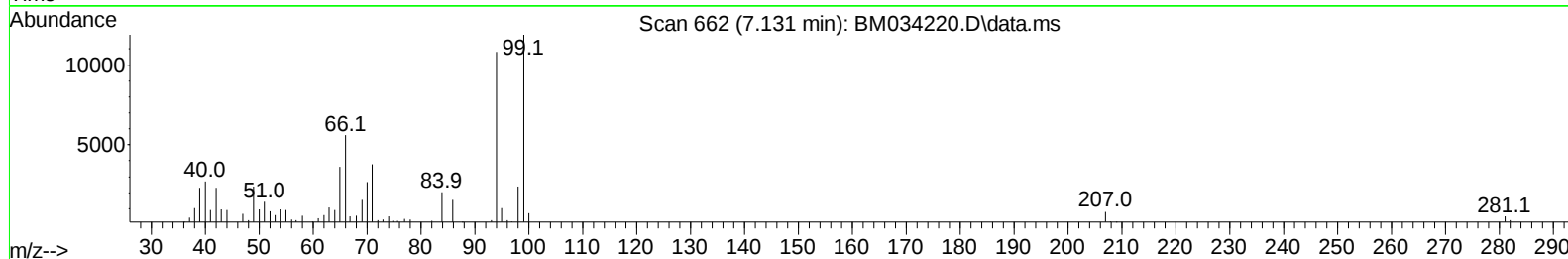
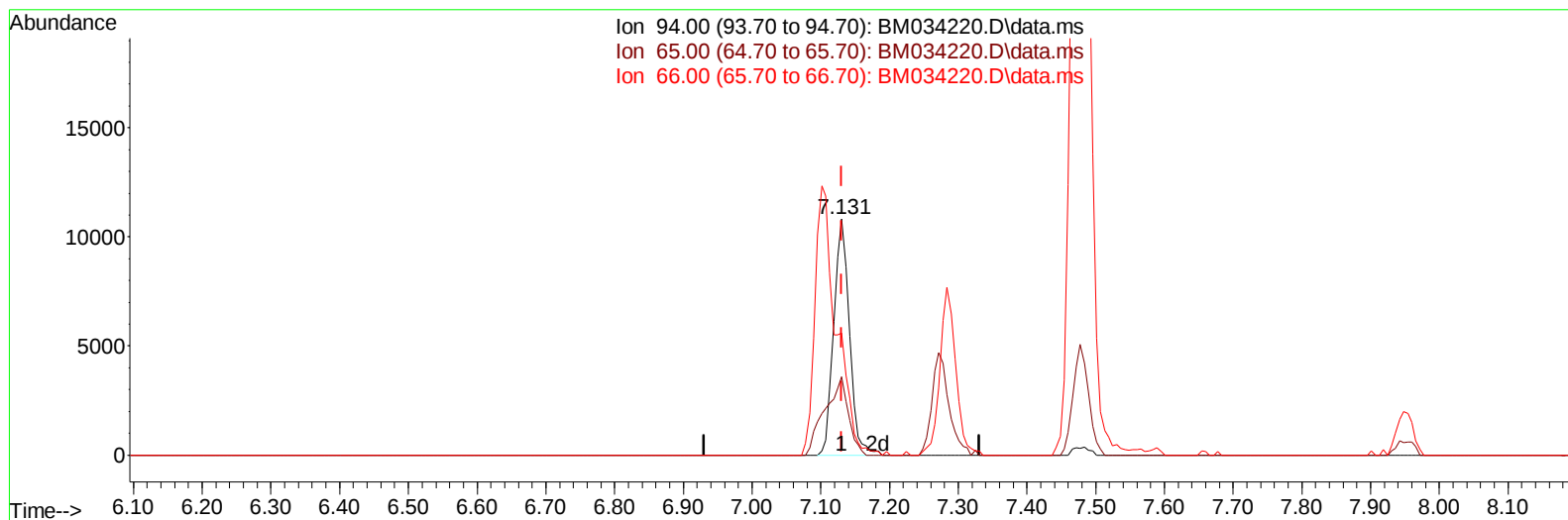
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TIC: BM034220.D\data.ms

(8) Phenol

7.131min (+ 0.000) 1.73 ng/u1 m

response 17146

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	28.60	33.28
66.00	42.10	51.63#
0.00	0.00	0.00

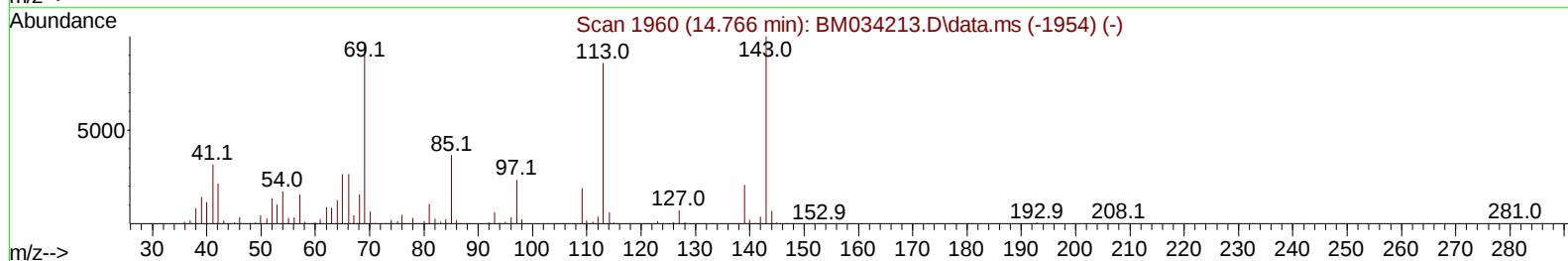
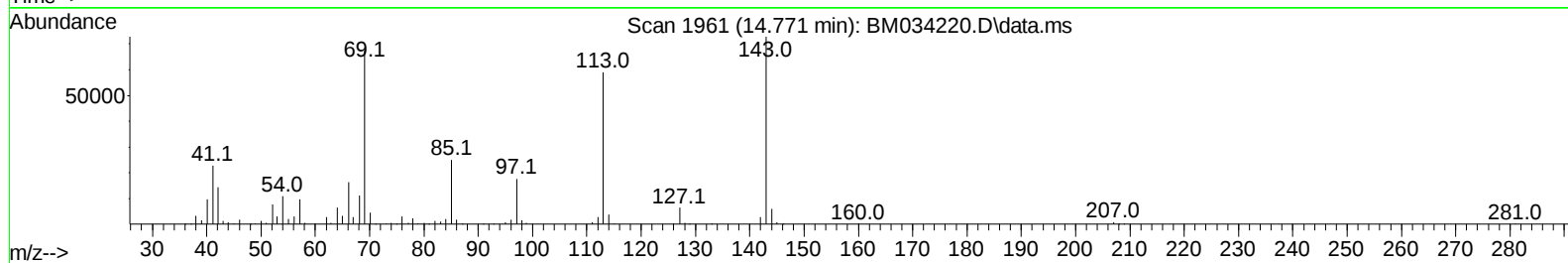
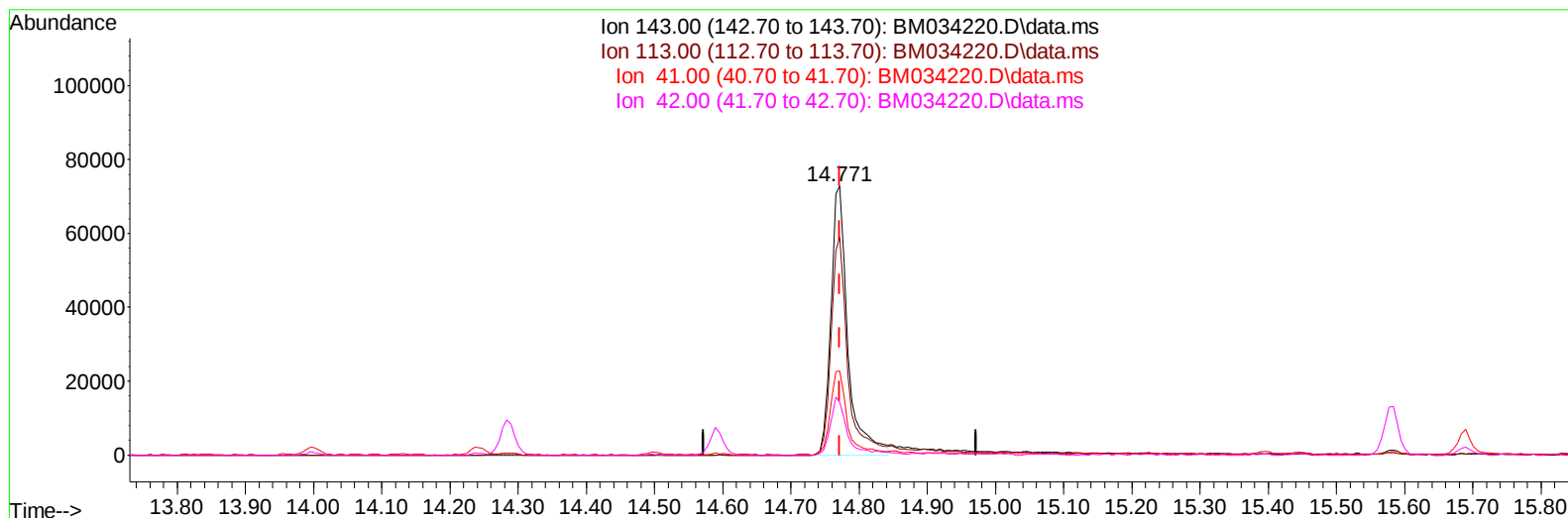
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Instrument :
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 ClientSampleId :
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TIC: BM034220.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.771min (+ 0.000) 28.36 ng/u1

response 125394

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	81.02
41.00	31.90	31.26
42.00	19.10	19.70

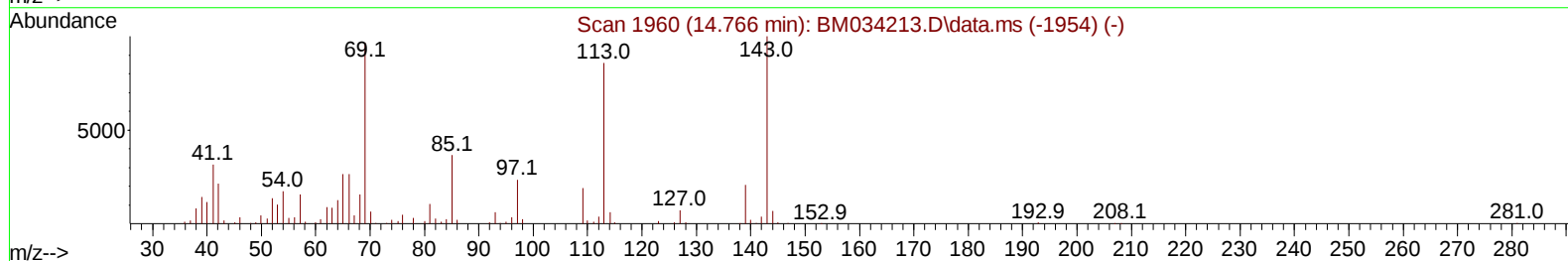
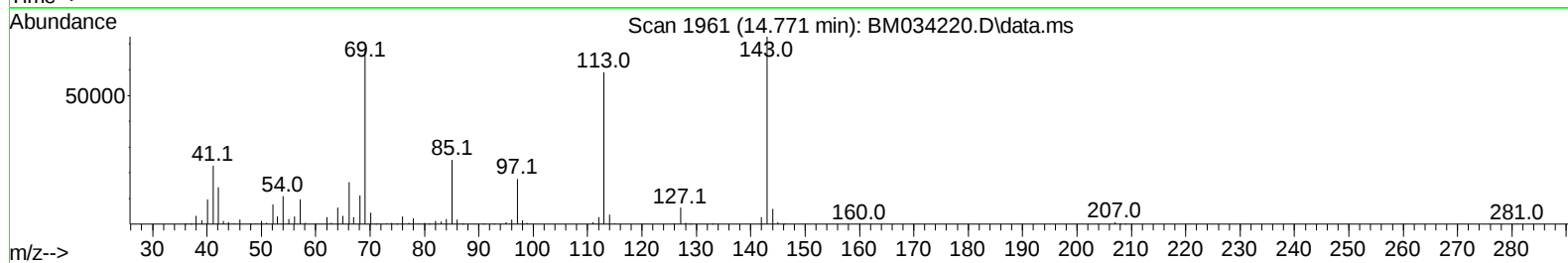
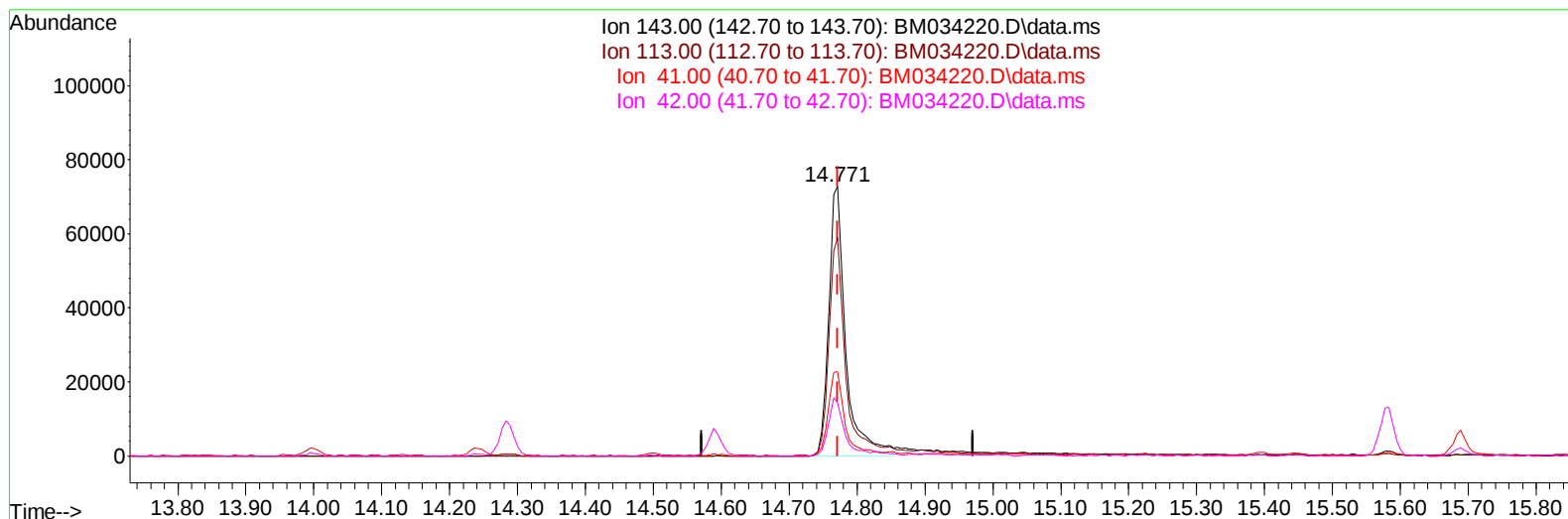
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 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
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TIC: BM034220.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.771min (+ 0.000) 28.76 ng/ul m

response 127142

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	81.02
41.00	31.90	31.26
42.00	19.10	19.70

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.954	152	149731	20.000	ng/ul	0.00
20) Naphthalene-d8	10.766	136	598016	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.595	164	390067	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.330	188	759281	20.000	ng/ul	0.00
79) Chrysene-d12	21.500	240	696359	20.000	ng/ul	0.00
88) Perylene-d12	23.924	264	780767	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	23623	6.887	ng/uL	0.00
4) Pyridine-d5	3.743	84	259151	31.733	ng/ul	0.00
7) Phenol-d5	7.101	99	356689	35.598	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.272	67	224667	38.220	ng/ul	0.00
11) 2-Chlorophenol-d4	7.478	132	333046	38.460	ng/ul	0.00
15) 4-Methylphenol-d8	8.654	113	316300	36.909	ng/ul	0.00
21) Nitrobenzene-d5	9.113	128	164883	38.439	ng/ul	0.00
24) 2-Nitrophenol-d4	9.842	143	188650	39.661	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.378	165	350700	36.778	ng/ul	0.00
31) 4-Chloroaniline-d4	10.895	131	428551	36.138	ng/ul	0.00
46) Dimethylphthalate-d6	14.001	166	1120488	39.479	ng/ul	0.00
49) Acenaphthylene-d8	14.283	160	1404042	38.949	ng/ul	0.00
54) 4-Nitrophenol-d4	14.771	143	127142m	28.759	ng/ul	0.00
60) Fluorene-d10	15.583	176	973421	39.180	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.689	200	148416	30.895	ng/ul	0.00
73) Anthracene-d10	17.430	188	1428786	39.841	ng/ul	0.00
81) Pyrene-d10	19.718	212	1667891	43.996	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.765	264	1511600	39.242	ng/ul	0.00
Target Compounds						
						Qvalue
8) Phenol	7.131	94	17146m	1.734	ng/ul	

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

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