

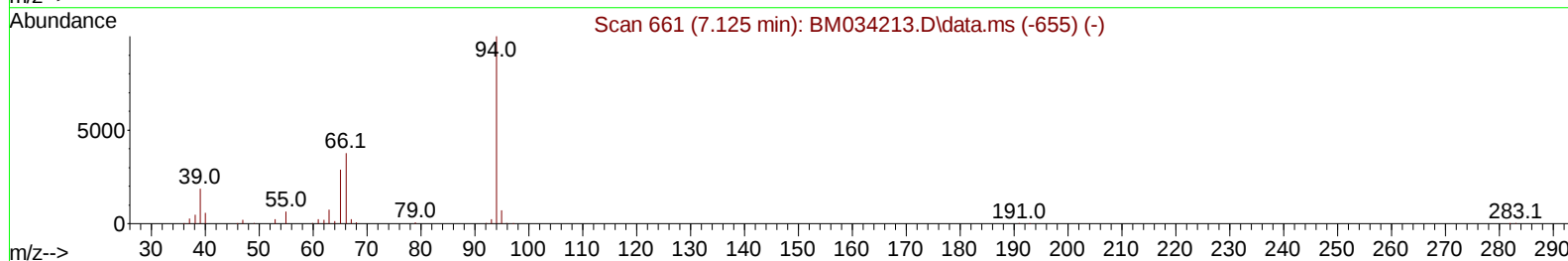
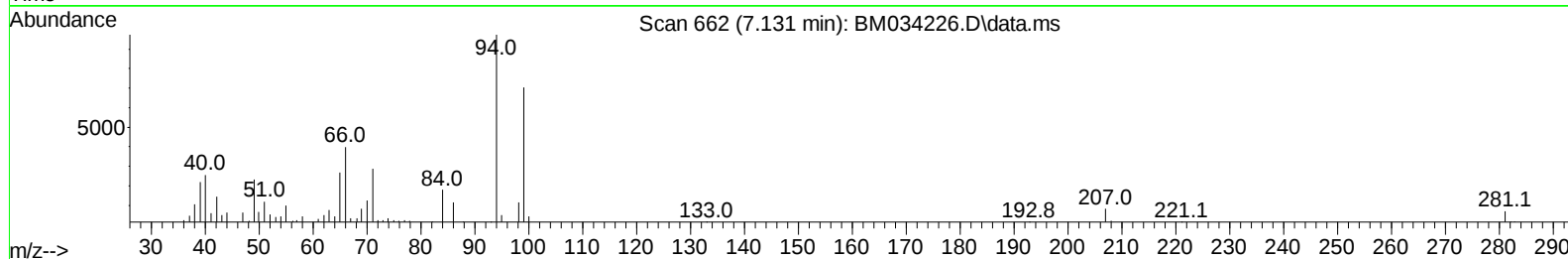
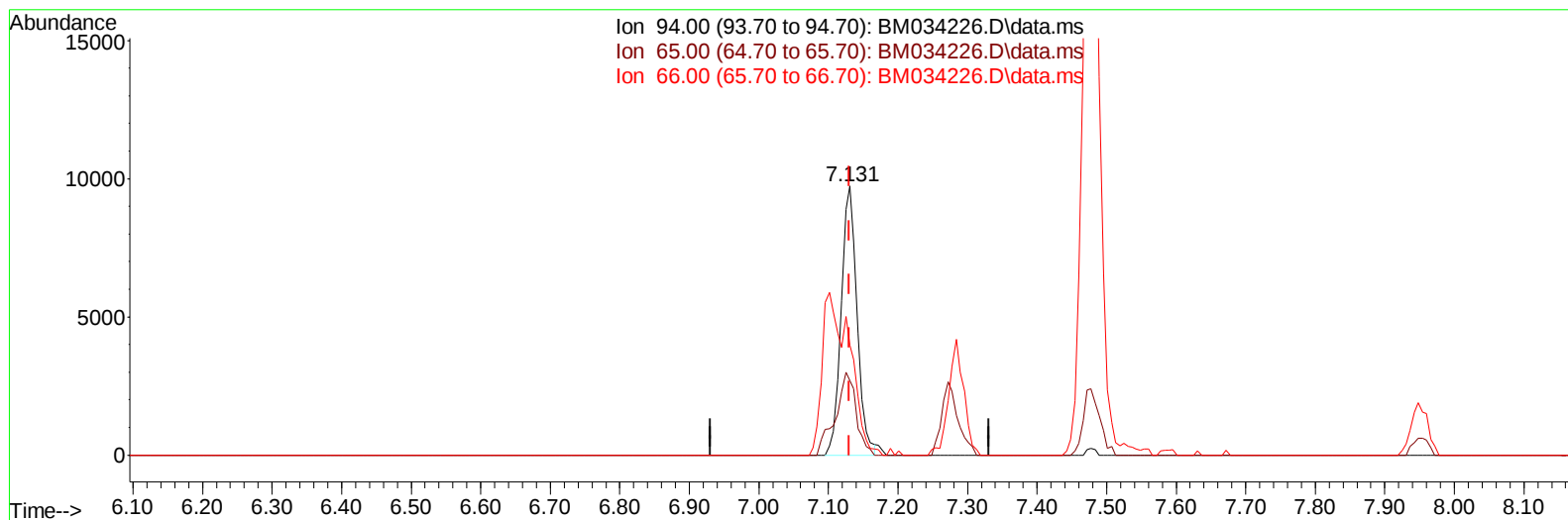
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031122\
 Data File : BM034226.D
 Acq On : 11 Mar 2022 17:24
 Operator : CG/JU
 Sample : N1816-13
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBSD8

Manual IntegrationsAPPROVED

Quant Time: Mar 11 23:41:31 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: BM034226.D\data.ms

(8) Phenol

7.131min (+ 0.000) 1.64 ng/u1

response 15499

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	28.60	27.66
66.00	42.10	41.02
0.00	0.00	0.00

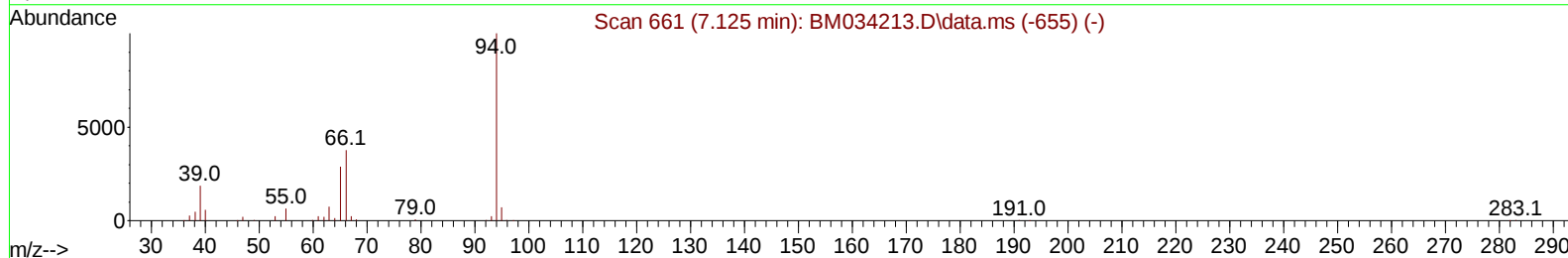
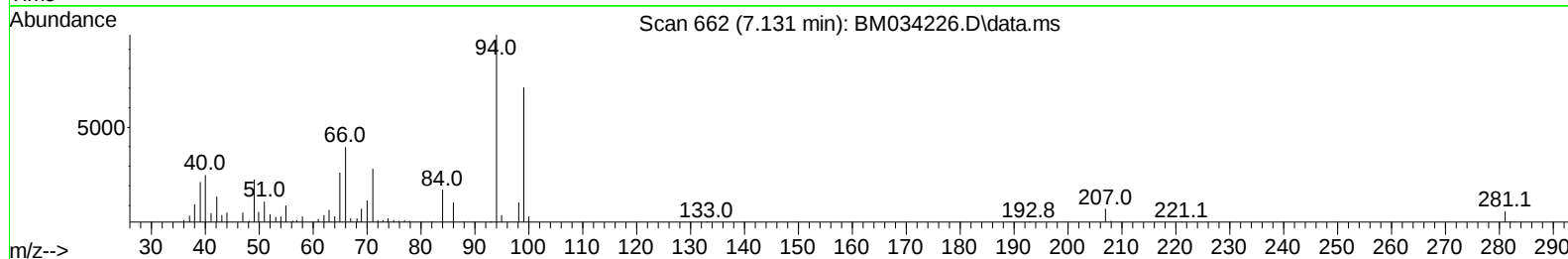
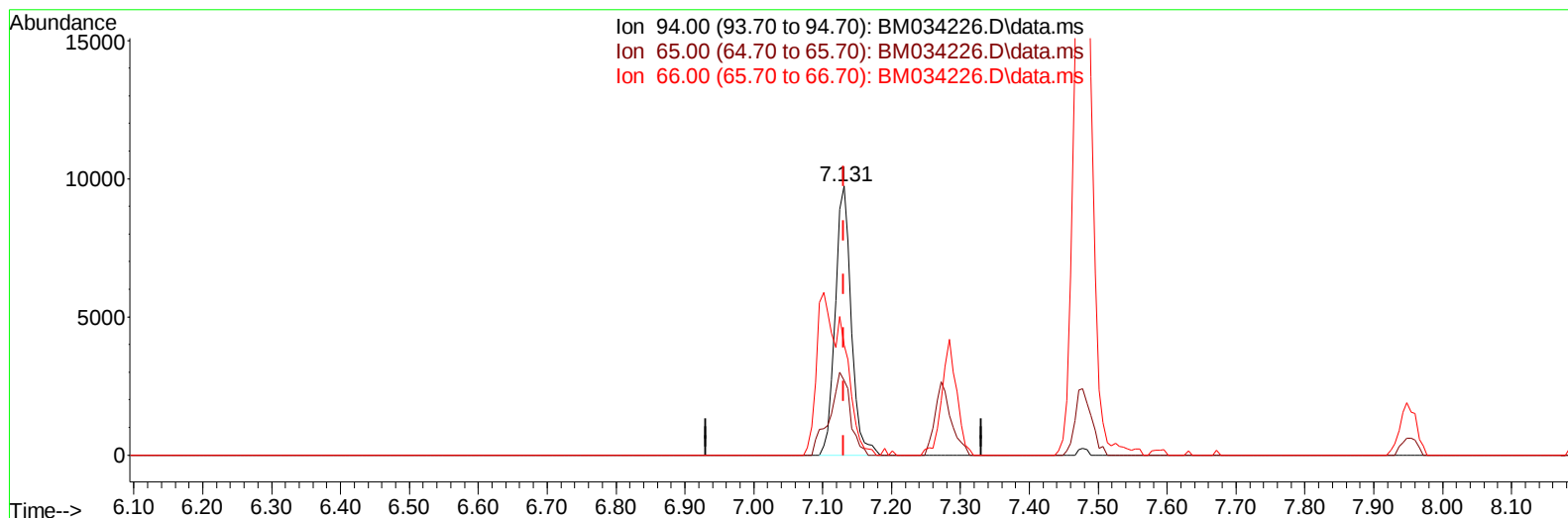
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TIC: BM034226.D\data.ms

(8) Phenol

7.131min (+ 0.000) 1.66 ng/u1 m

response 15682

Ion	Exp%	Act%
94.00	100.00	100.00
65.00	28.60	27.66
66.00	42.10	41.02
0.00	0.00	0.00

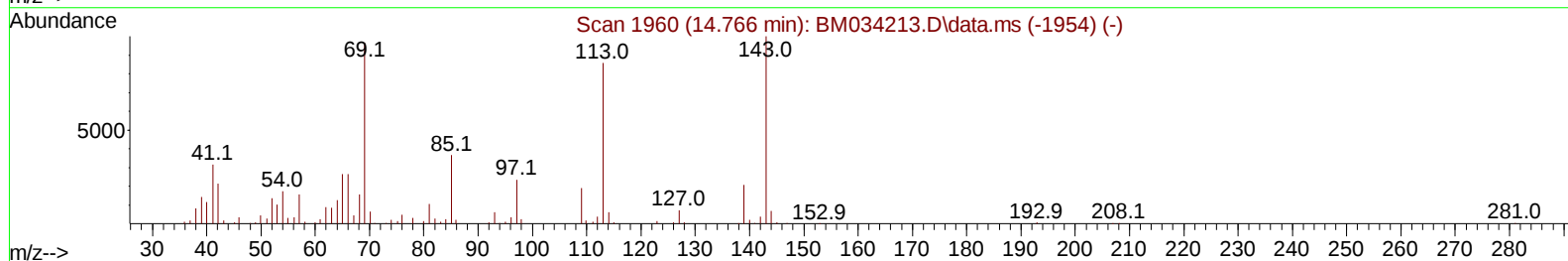
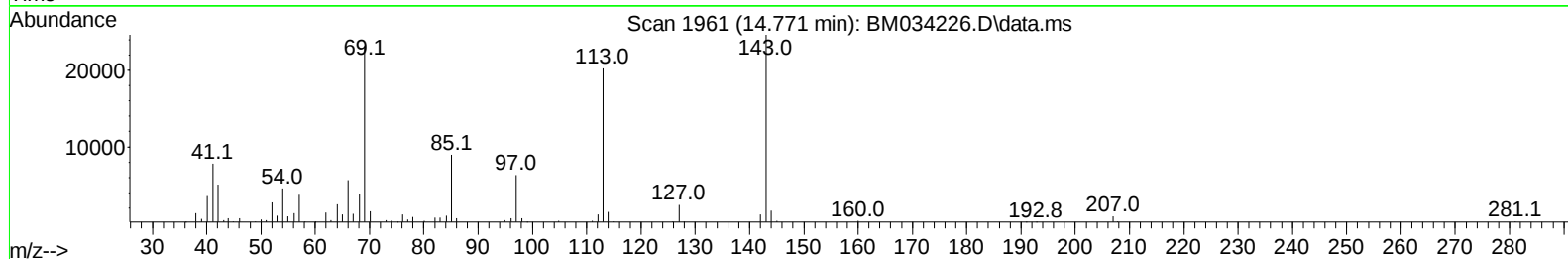
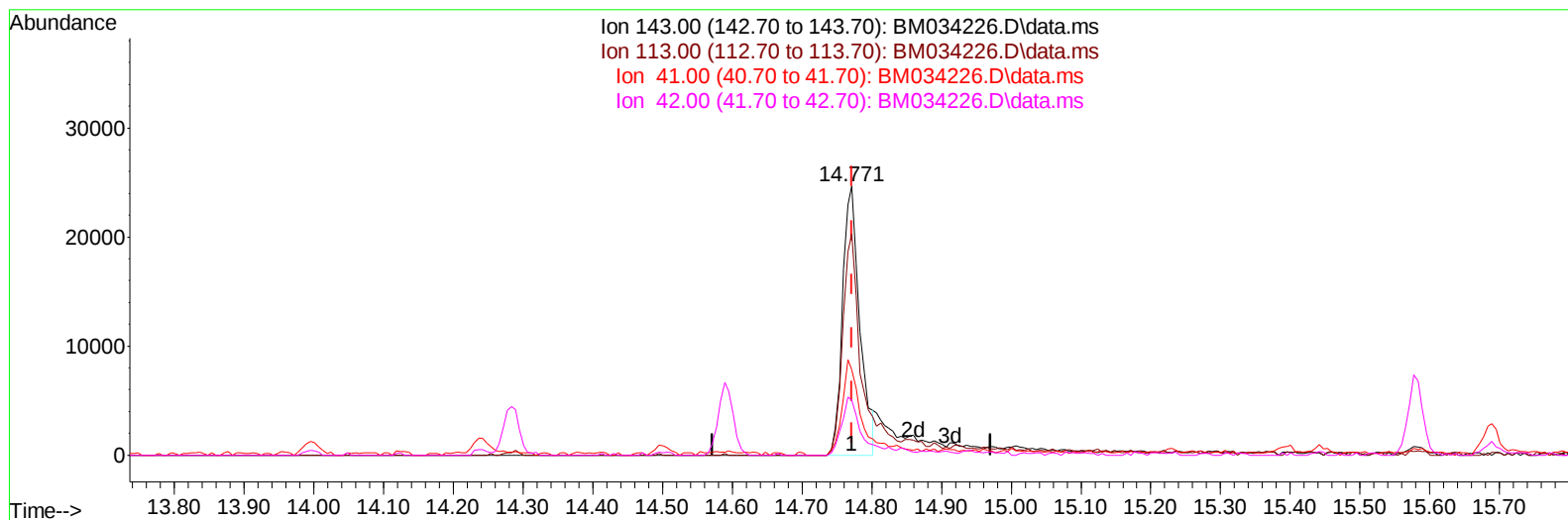
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 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
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TIC: BM034226.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.771min (+ 0.000) 9.85 ng/u1

response 42046

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	82.10
41.00	31.90	31.73
42.00	19.10	20.41

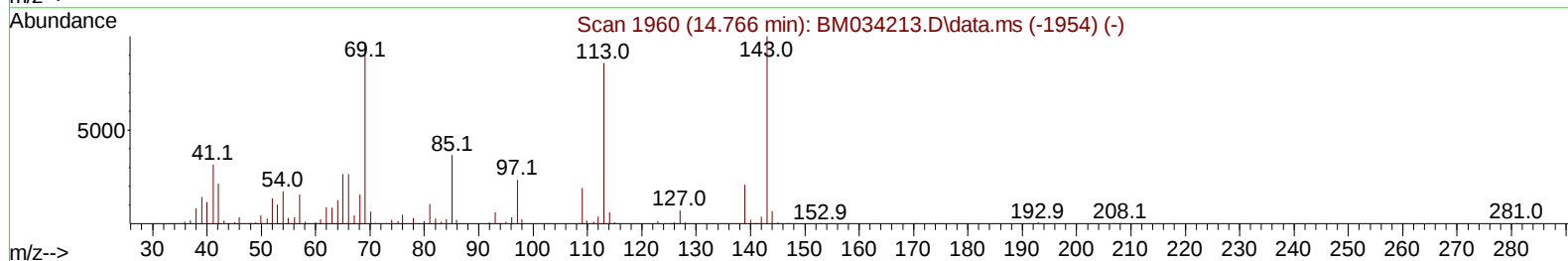
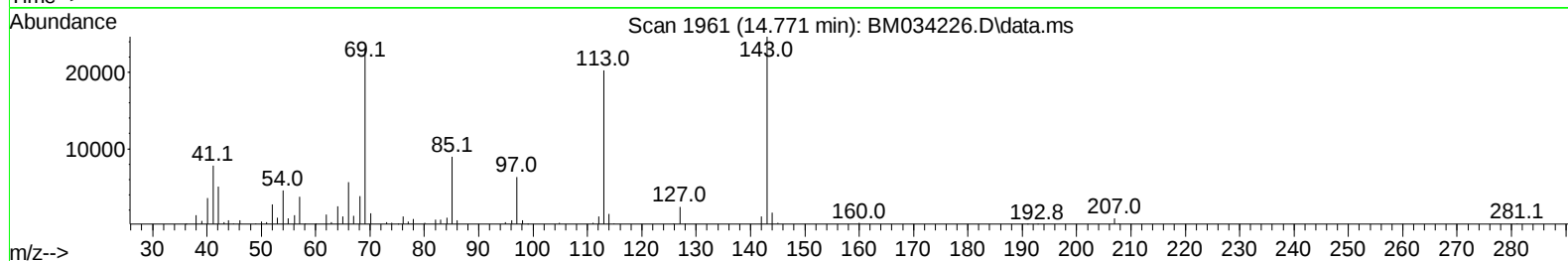
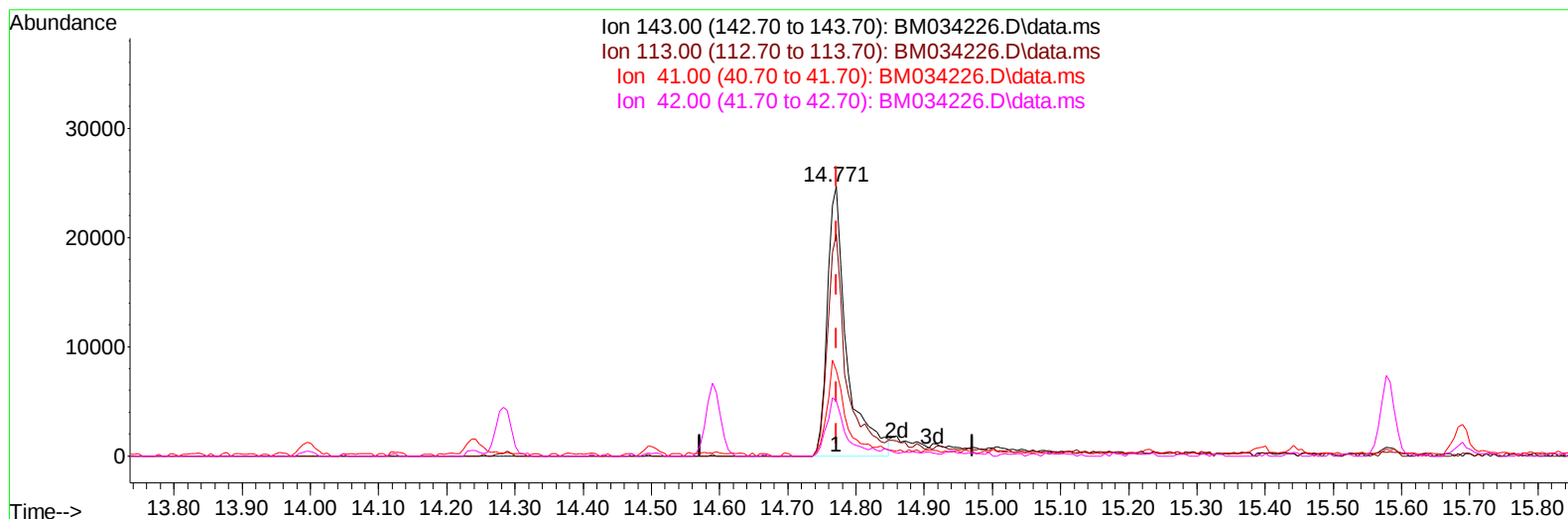
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 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
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TIC: BM034226.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.771min (+ 0.000) 11.45 ng/ul m

response 48856

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	82.10
41.00	31.90	31.73
42.00	19.10	20.41

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.954	152	142794	20.000	ng/ul	0.00
20) Naphthalene-d8	10.766	136	566724	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.589	164	376456	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.330	188	745761	20.000	ng/ul	0.00
79) Chrysene-d12	21.500	240	701961	20.000	ng/ul	0.00
88) Perylene-d12	23.924	264	764700	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	12374	3.783	ng/uL	0.00
4) Pyridine-d5	3.743	84	110539	14.193	ng/ul	0.00
7) Phenol-d5	7.101	99	176242	18.443	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.272	67	116123	20.714	ng/ul	0.00
11) 2-Chlorophenol-d4	7.478	132	169629	20.540	ng/ul	0.00
15) 4-Methylphenol-d8	8.654	113	152324	18.638	ng/ul	0.00
21) Nitrobenzene-d5	9.113	128	82088	20.193	ng/ul	0.00
24) 2-Nitrophenol-d4	9.842	143	86107	19.102	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.378	165	170673	18.887	ng/ul	0.00
31) 4-Chloroaniline-d4	10.895	131	204373	18.186	ng/ul	0.00
46) Dimethylphthalate-d6	13.995	166	578217	21.109	ng/ul	0.00
49) Acenaphthylene-d8	14.283	160	711946	20.464	ng/ul	0.00
54) 4-Nitrophenol-d4	14.771	143	48856m	11.451	ng/ul	0.00
60) Fluorene-d10	15.577	176	510742	21.300	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.689	200	67869	14.384	ng/ul	0.00
73) Anthracene-d10	17.430	188	762634	21.651	ng/ul	0.00
81) Pyrene-d10	19.718	212	905678	23.699	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.765	264	796843	21.121	ng/ul	0.00
Target Compounds						
8) Phenol	7.131	94	15682m	1.663	ng/ul	Qvalue

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

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