

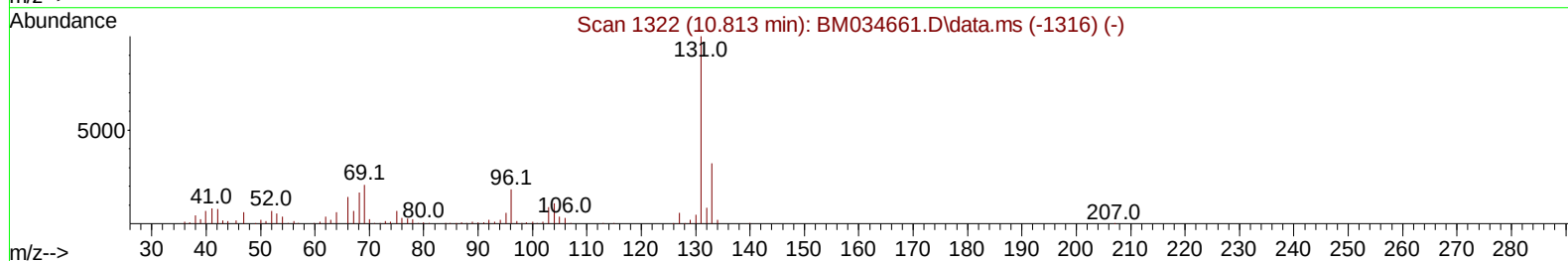
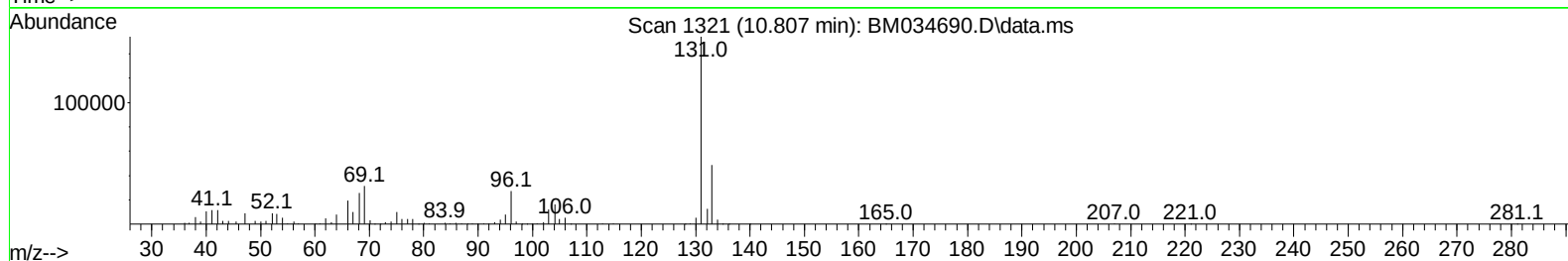
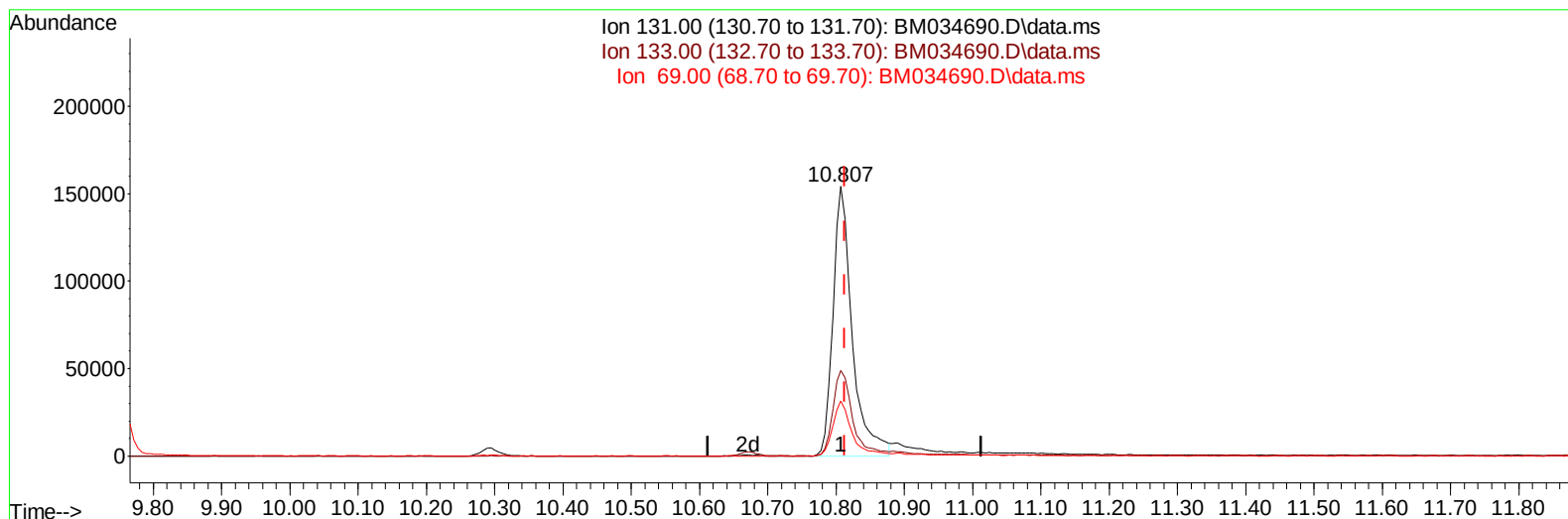
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034690.D
Acq On : 15 Apr 2022 04:51
Operator : CG/JU
Sample : N2316-11
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNR1

Manual IntegrationsAPPROVED

Quant Time: Apr 15 06:09:05 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 14 23:42:00 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/15/2022
Supervised By :Yogesh Patel 04/23/2022



TIC: BM034690.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.807min (-0.006) 30.04 ng/ul

response 303010

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.00	31.63
69.00	16.80	20.36#
0.00	0.00	0.00

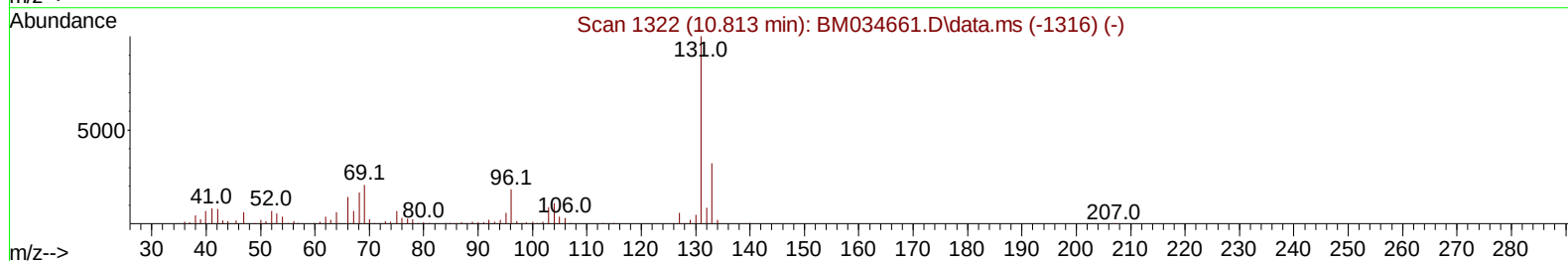
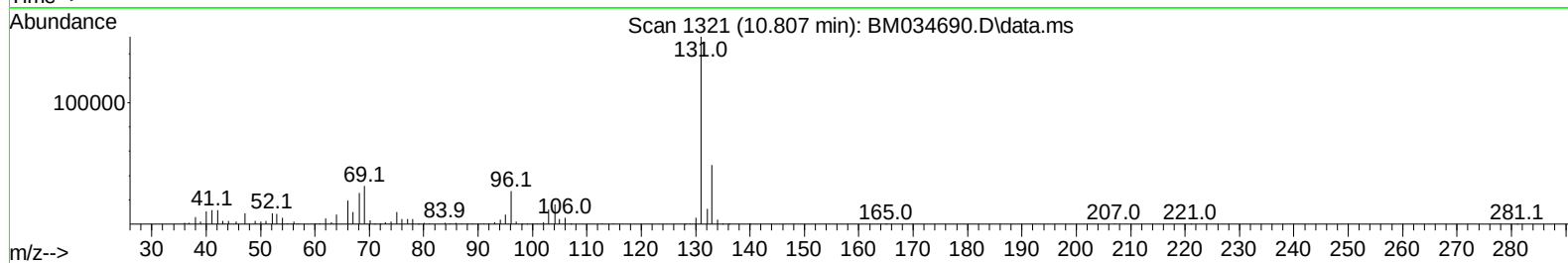
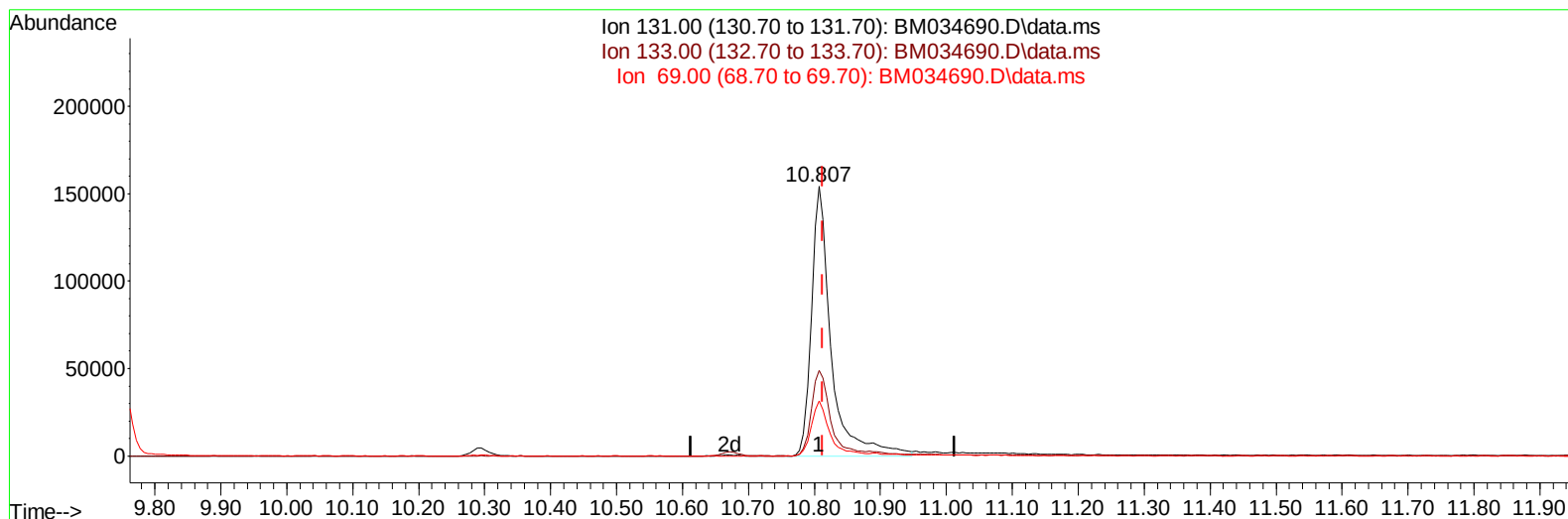
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TIC: BM034690.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.807min (-0.006) 32.00 ng/ul m

response 322767

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.00	31.63
69.00	16.80	20.36#
0.00	0.00	0.00

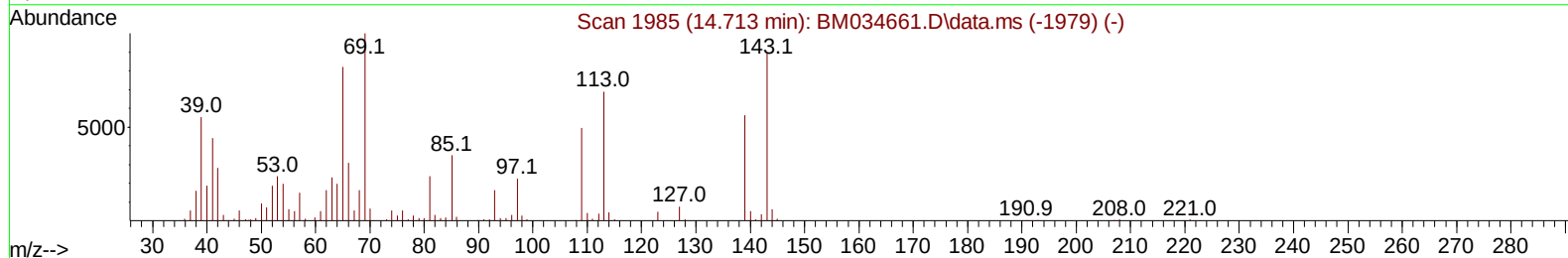
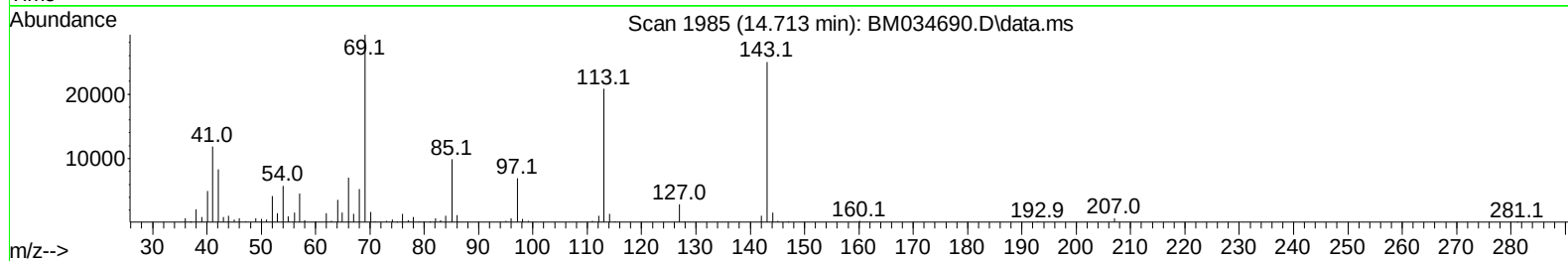
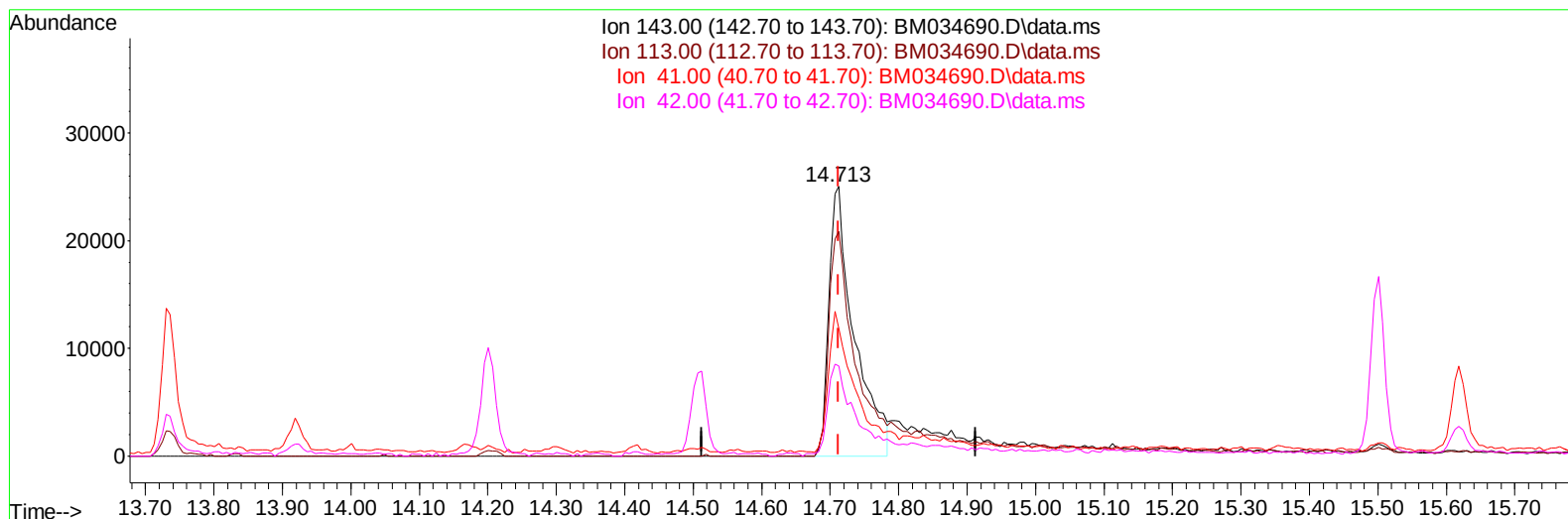
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TIC: BM034690.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.713min (-0.000) 19.47 ng/ul

response 64335

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	83.25
41.00	31.90	47.31#
42.00	19.10	33.41#

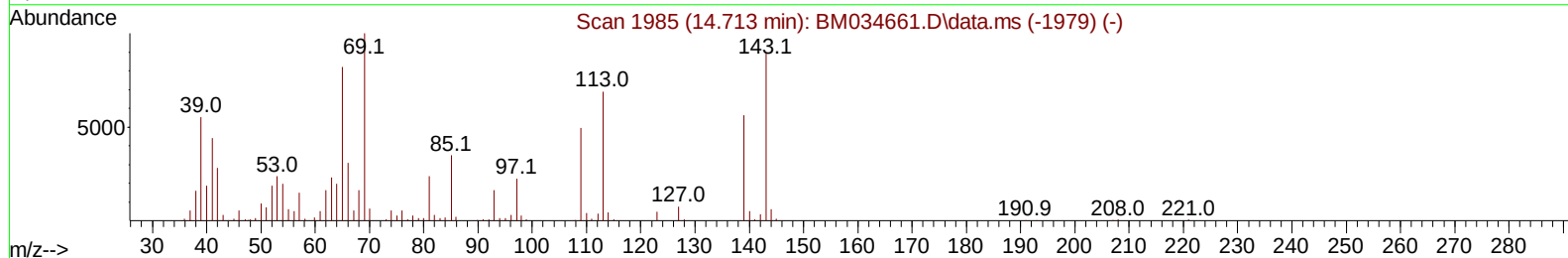
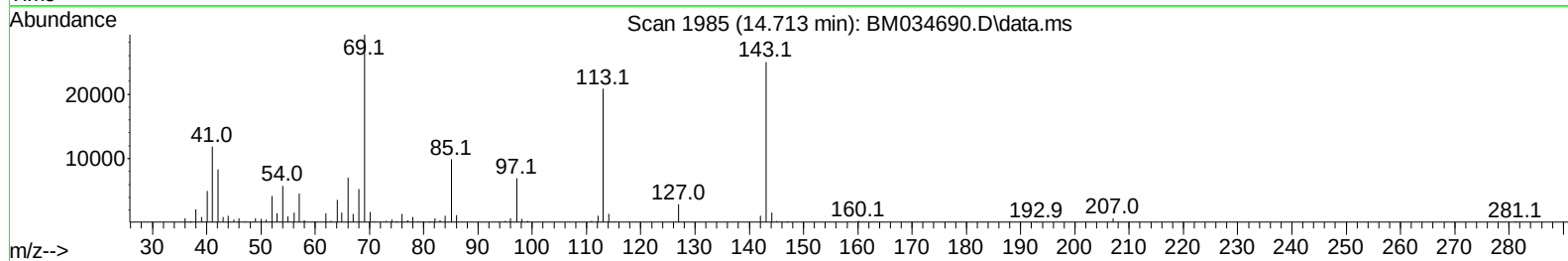
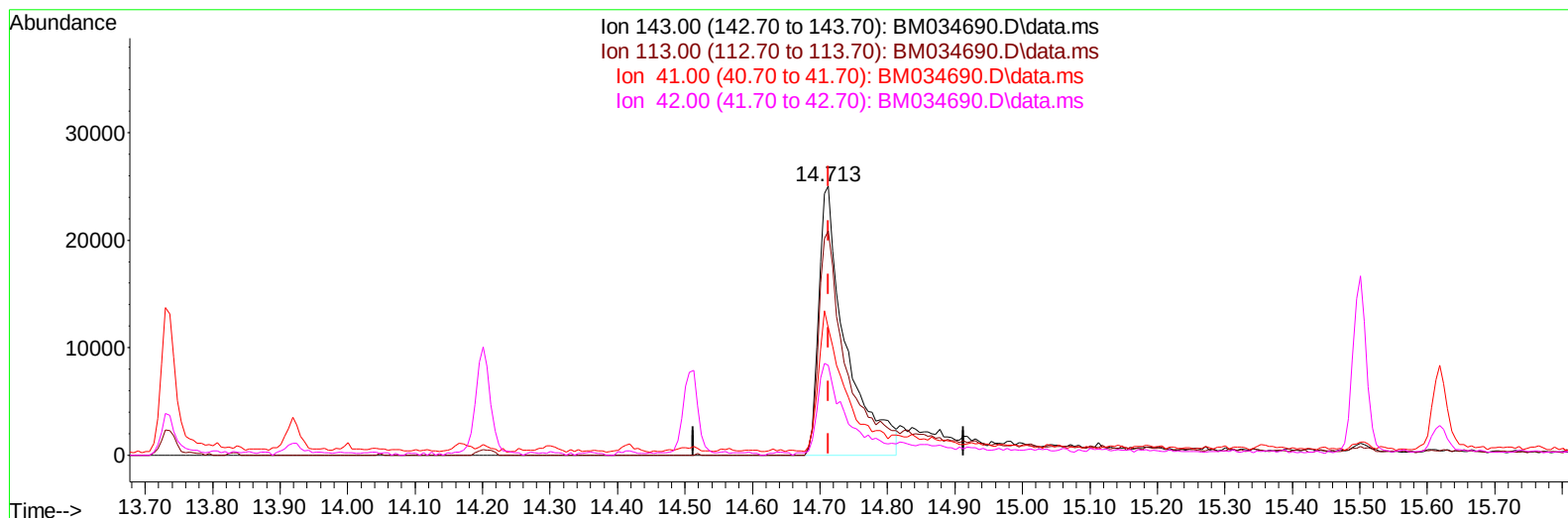
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 Sample : N2316-11
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BM034690.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.713min (-0.000) 21.06 ng/ul m

response 69590

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	83.25
41.00	31.90	47.31#
42.00	19.10	33.41#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.860	152	121506	20.000	ng/ul	0.00
20) Naphthalene-d8	10.672	136	507451	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.507	164	299979	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.254	188	583067	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.430	240	546668	20.000	ng/ul	# 0.00
88) Perylene-d12	23.800	264	629579	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.243	96	16405	4.815	ng/uL	0.00
4) Pyridine-d5	3.666	84	230000	27.970	ng/ul	0.00
7) Phenol-d5	7.025	99	291666	28.840	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.190	67	213219	31.640	ng/ul	0.00
11) 2-Chlorophenol-d4	7.390	132	241808	30.930	ng/ul	0.00
15) 4-Methylphenol-d8	8.572	113	263008	33.293	ng/ul	-0.01
21) Nitrobenzene-d5	9.025	128	125980	33.779	ng/ul	0.00
24) 2-Nitrophenol-d4	9.748	143	135301	34.846	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.289	165	227155	31.068	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.807	131	322767m	31.998	ng/ul	0.00
46) Dimethylphthalate-d6	13.919	166	811907	37.457	ng/ul	0.00
49) Acenaphthylene-d8	14.201	160	976523	35.051	ng/ul	0.00
54) 4-Nitrophenol-d4	14.713	143	69590m	21.055	ng/ul	0.00
60) Fluorene-d10	15.501	176	700827	37.414	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.618	200	103599	29.105	ng/ul	0.00
73) Anthracene-d10	17.348	188	1030127	37.057	ng/ul	0.00
81) Pyrene-d10	19.642	212	1191146	39.922	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.647	264	1115389	35.220	ng/ul	-0.01

Target Compounds Qvalue

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

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