

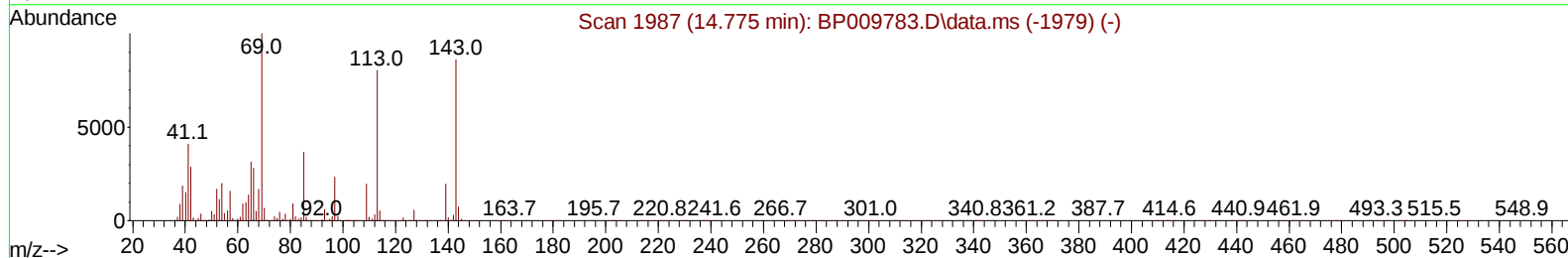
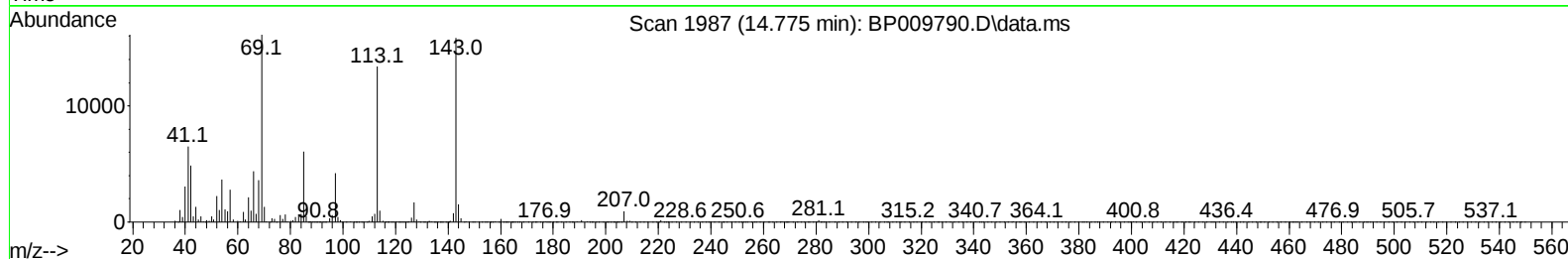
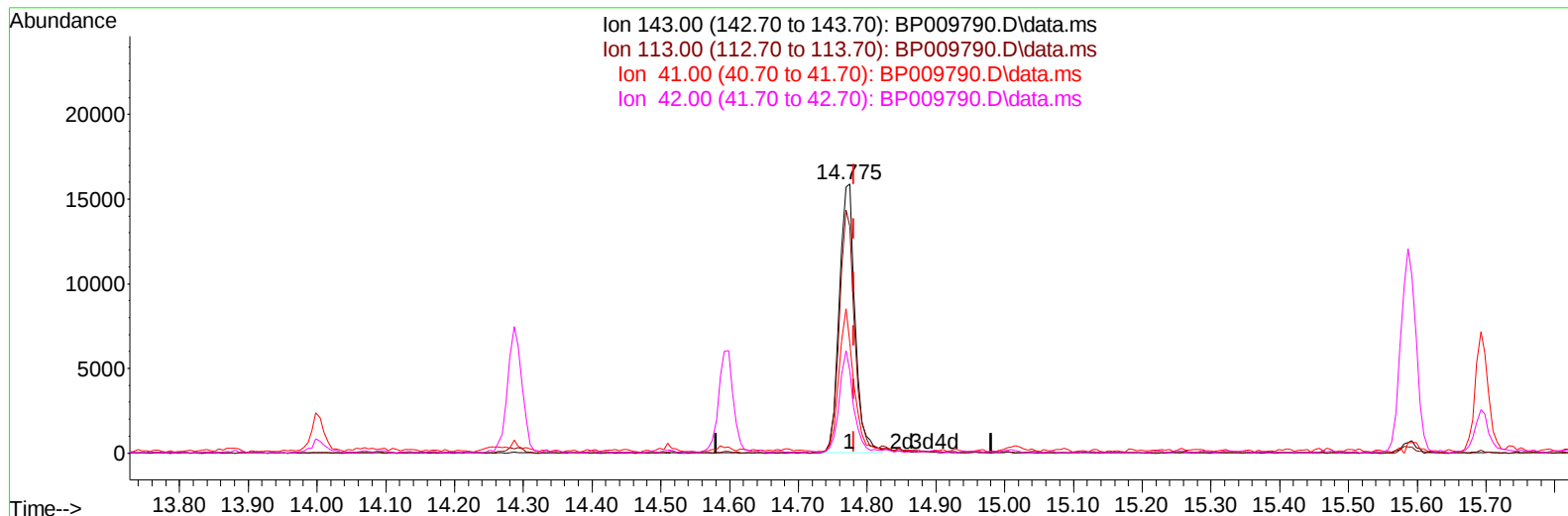
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009790.D
 Acq On : 12 Apr 2022 19:16
 Operator : CG/JU
 Sample : N2342-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J64

Manual IntegrationsAPPROVED

Quant Time: Apr 13 02:24:52 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 13 02:15:22 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/13/2022
 Supervised By :Yogesh Patel 04/13/2022



TIC: BP009790.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.006) 7.21 ng/u1

response 25520

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	84.27#
41.00	23.20	40.95#
42.00	18.00	30.63#

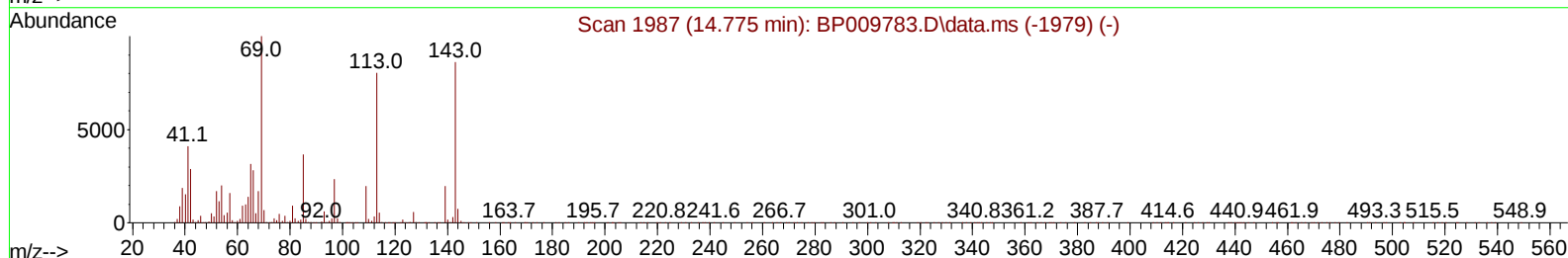
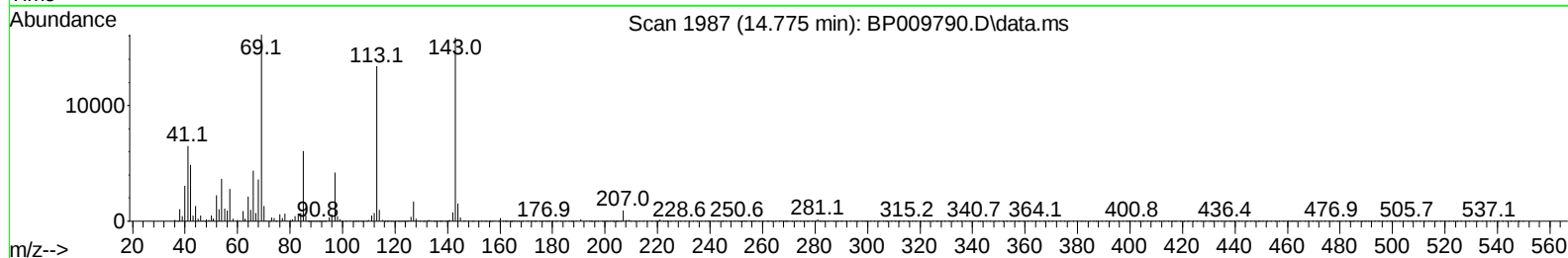
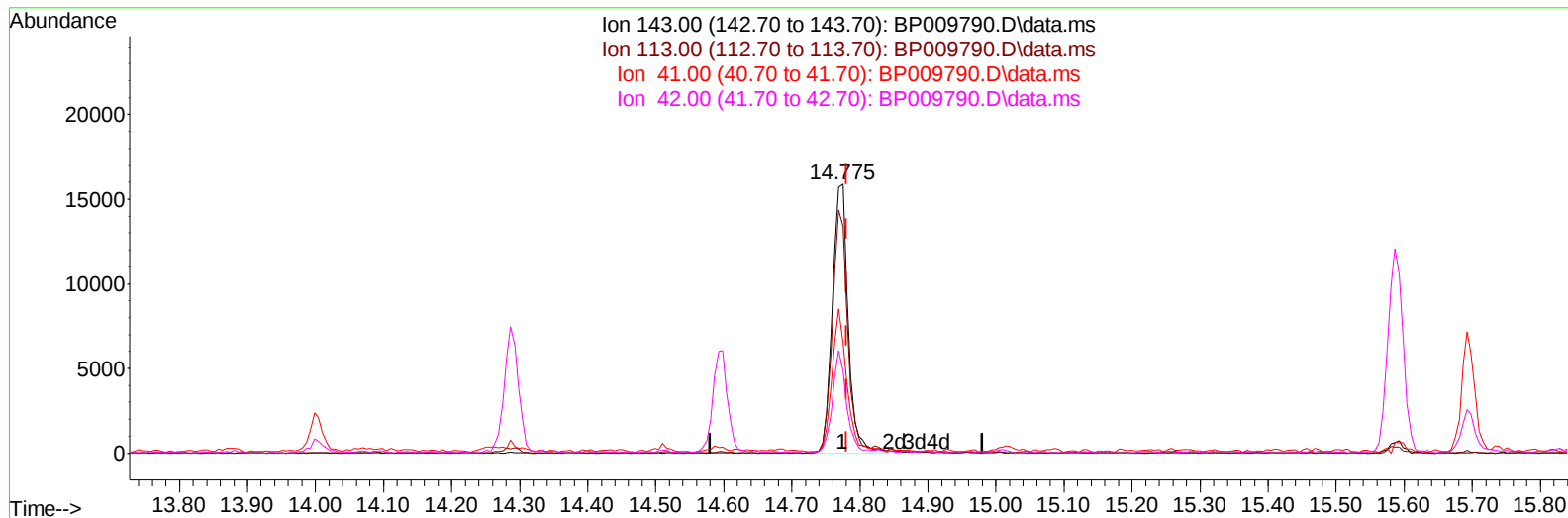
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TIC: BP009790.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.006) 7.28 ng/ul m

response 25771

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	84.27#
41.00	23.20	40.95#
42.00	18.00	30.63#

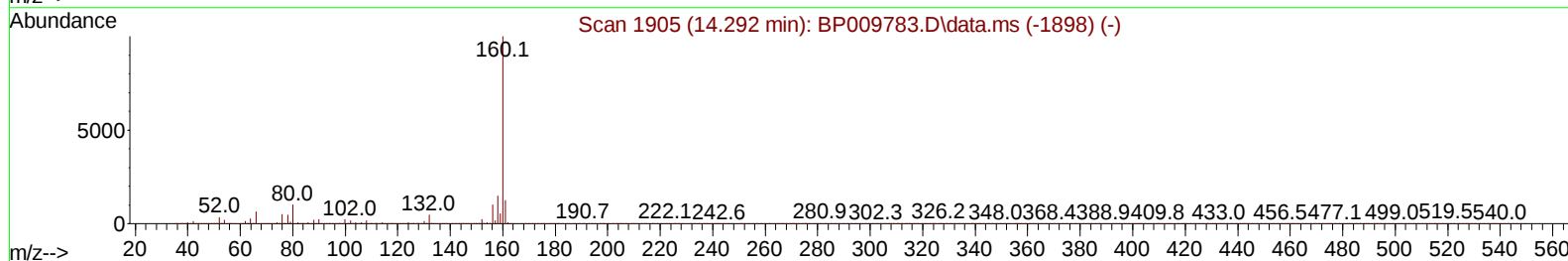
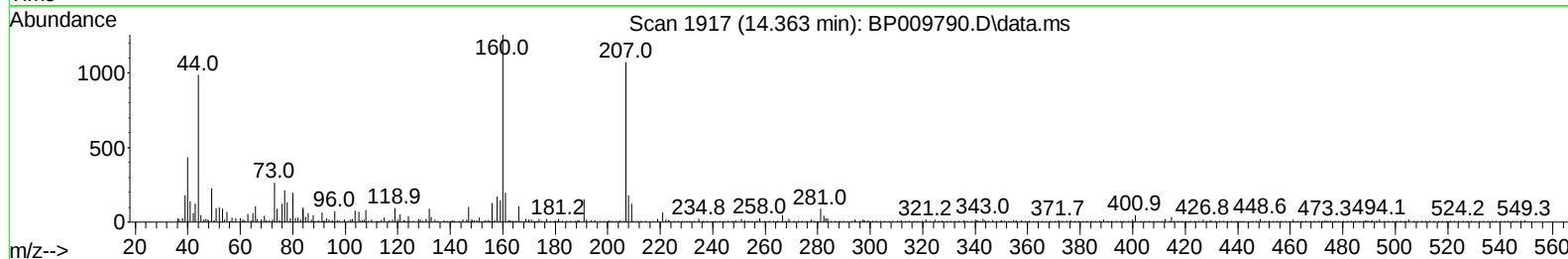
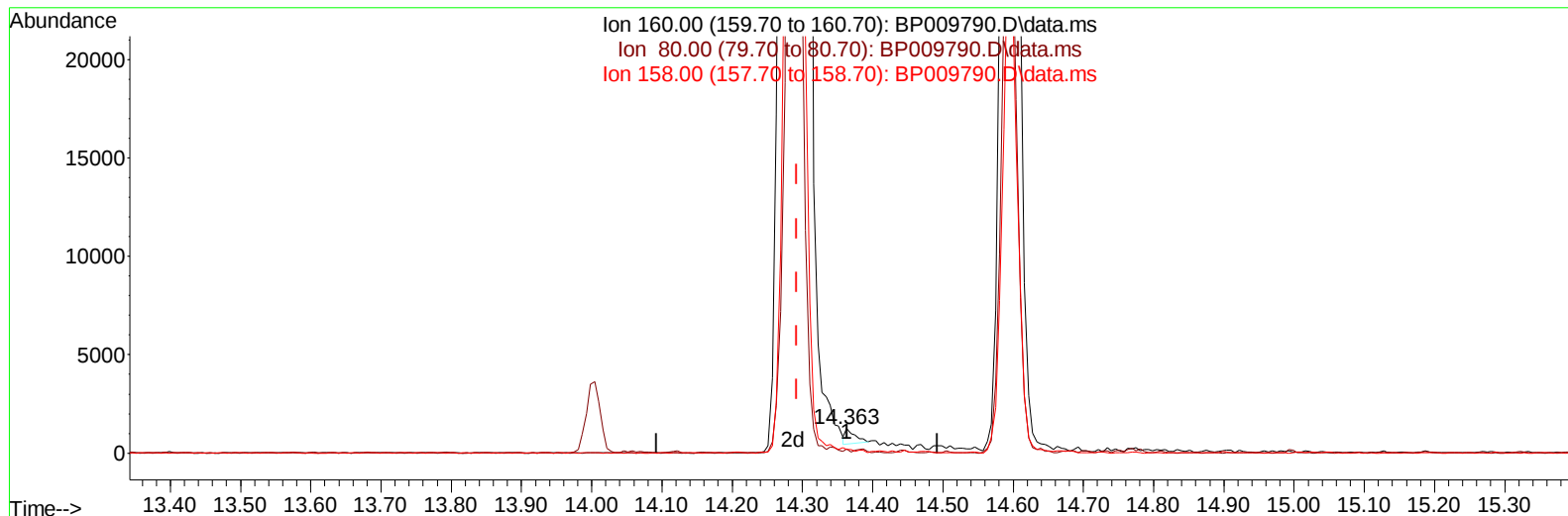
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TIC: BP009790.D\data.ms

(49) Acenaphthylene-d8 (S)

14.363min (+ 0.071) 0.03 ng/u1

response 764

Ion	Exp%	Act%
160.00	100.00	100.00
80.00	17.00	15.76
158.00	13.10	13.61
0.00	0.00	0.00

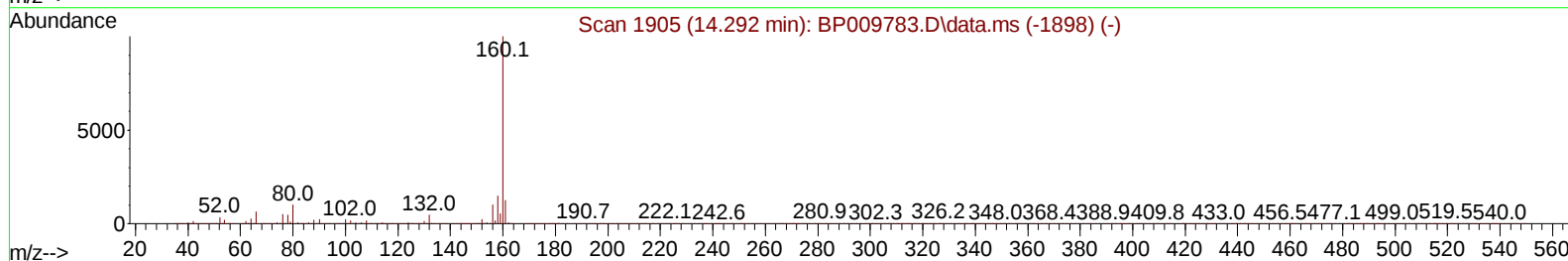
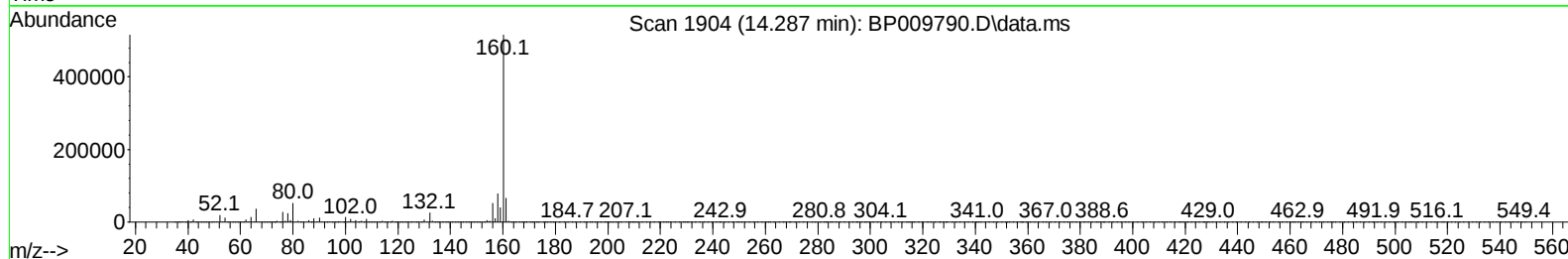
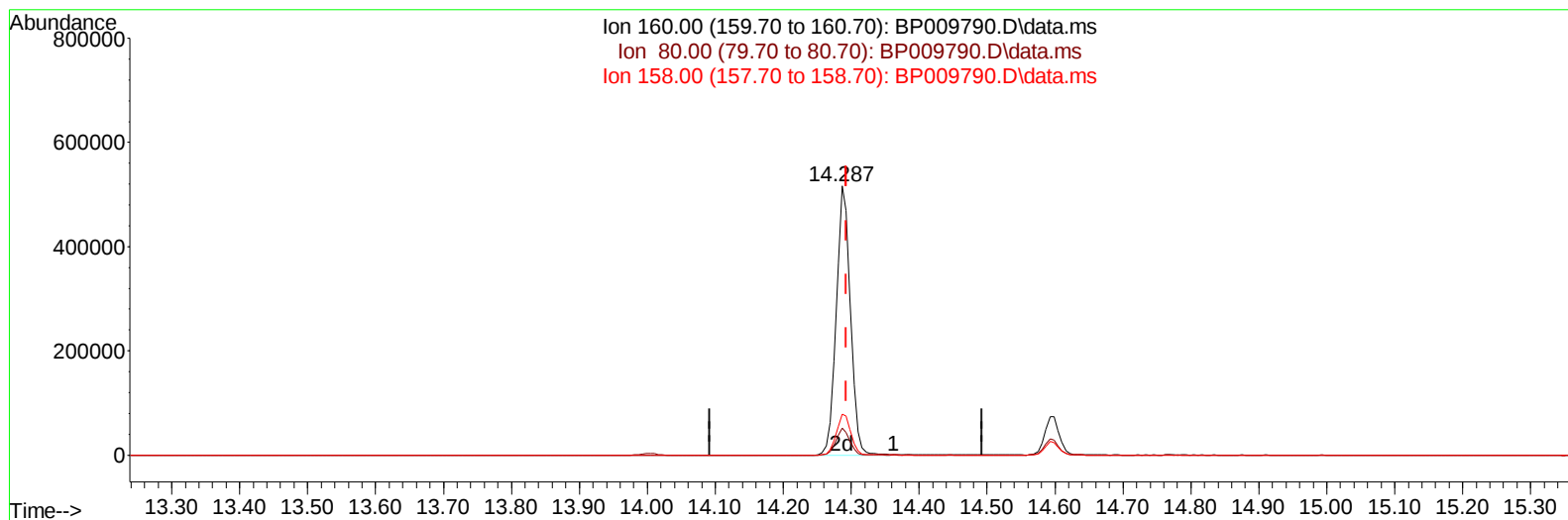
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Data File : BP009790.D
Acq On : 12 Apr 2022 19:16
Operator : CG/JU
Sample : N2342-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP009790.D\data.ms

(49) Acenaphthylene-d8 (S)

14.287min (-0.006) 32.52 ng/ul m

response 752286

Ion	Exp%	Act%
160.00	100.00	100.00
80.00	17.00	10.13#
158.00	13.10	15.13
0.00	0.00	0.00

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 Misc :
 ALS Vial : 9 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	76767	20.000	ng/ul	0.00
20) Naphthalene-d8	10.775	136	342290	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.598	164	248557	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	565634	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.427	240	644891	20.000	ng/ul	# 0.00
88) Perylene-d12	23.892	264	691911	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	8099	4.708	ng/uL	0.00
4) Pyridine-d5	3.805	84	33634	7.092	ng/ul	0.00
7) Phenol-d5	7.116	99	52093	7.729	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	140715	35.070	ng/ul	0.00
11) 2-Chlorophenol-d4	7.493	132	134400	26.437	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	100230	17.815	ng/ul	0.00
21) Nitrobenzene-d5	9.122	128	90308	36.586	ng/ul	0.00
24) 2-Nitrophenol-d4	9.852	143	89677	33.138	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	162051	28.617	ng/ul	0.00
31) 4-Chloroaniline-d4	10.904	131	229924	28.507	ng/ul	0.00
46) Dimethylphthalate-d6	14.004	166	694277	35.559	ng/ul	0.00
49) Acenaphthylene-d8	14.287	160	752286m	32.523	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	25771m	7.278	ng/ul	0.00
60) Fluorene-d10	15.587	176	614095	37.389	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	96544	28.812	ng/ul	0.00
73) Anthracene-d10	17.445	188	1014971	37.478	ng/ul	0.00
81) Pyrene-d10	19.675	212	1285707	36.504	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.733	264	1360216	37.750	ng/ul	0.00

Target Compounds Qvalue

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

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