

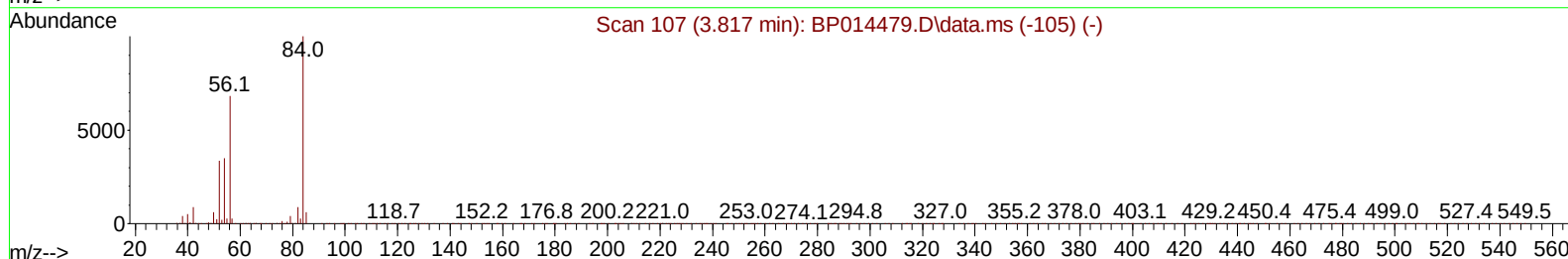
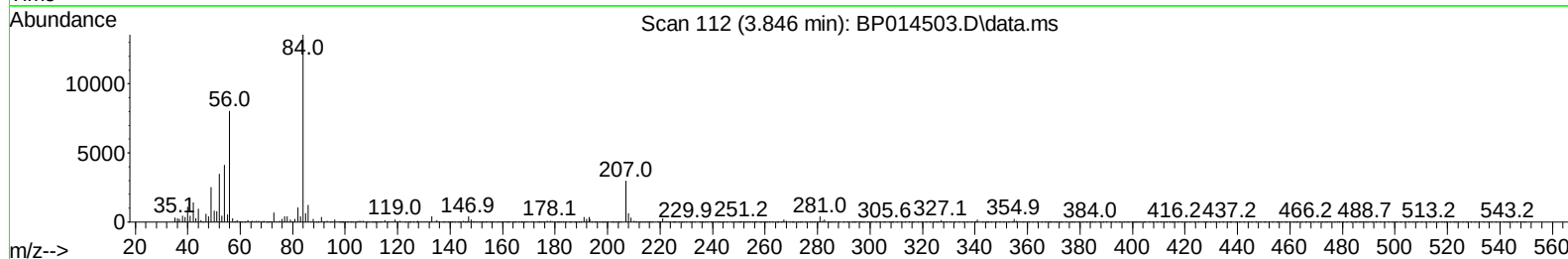
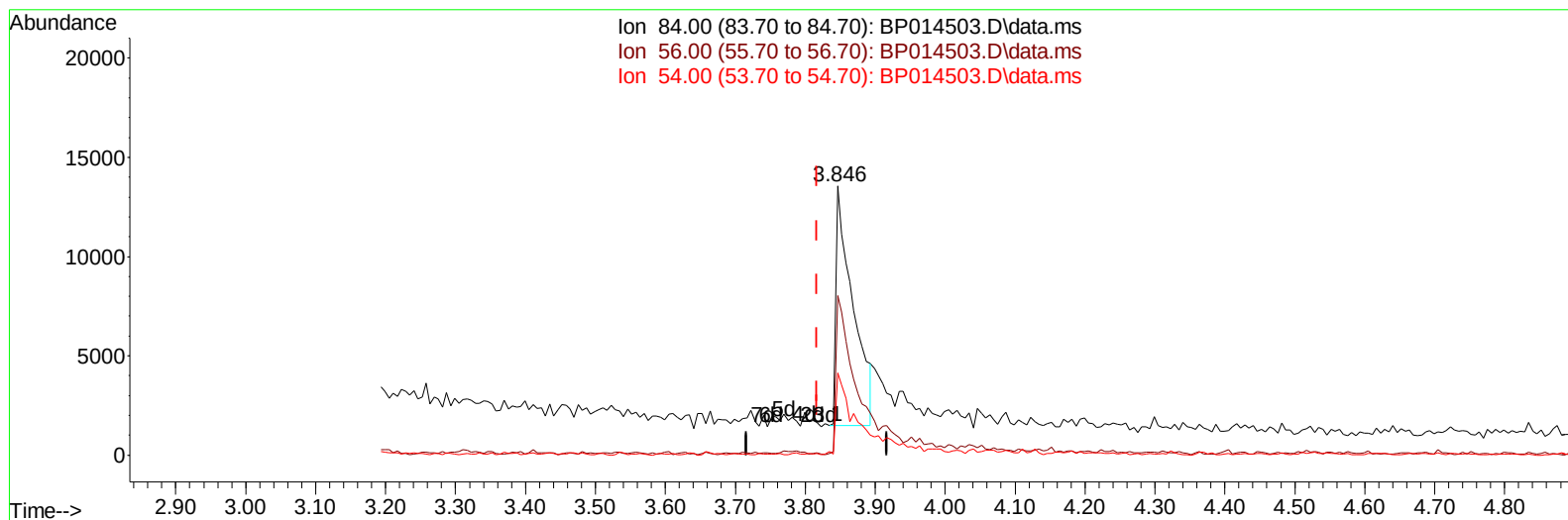
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014503.D
Acq On : 04 Apr 2023 00:17
Operator : CG/JU
Sample : 02158-07
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PMW-6(20230329)

Manual IntegrationsAPPROVED

Quant Time: Apr 04 01:40:48 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Apr 04 01:34:36 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/04/2023
Supervised By :Jagrut Upadhyay 04/04/2023



TIC: BP014503.D\data.ms

(4) Pyridine-d5 (S)

3.846min (+ 0.029) 2.12 ng/u1

response 20445

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	59.29
54.00	32.60	30.56
0.00	0.00	0.00

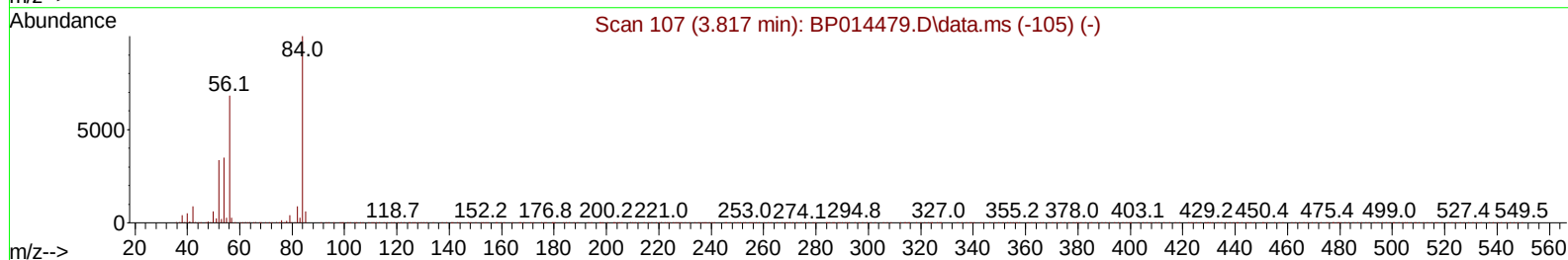
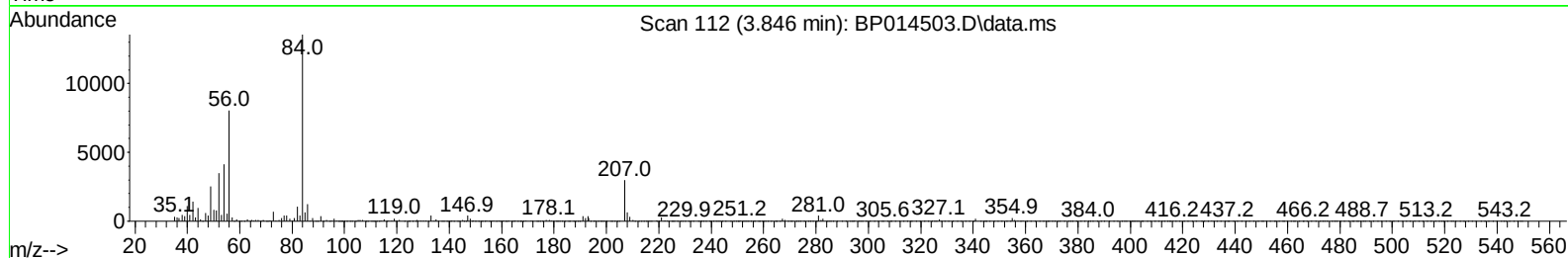
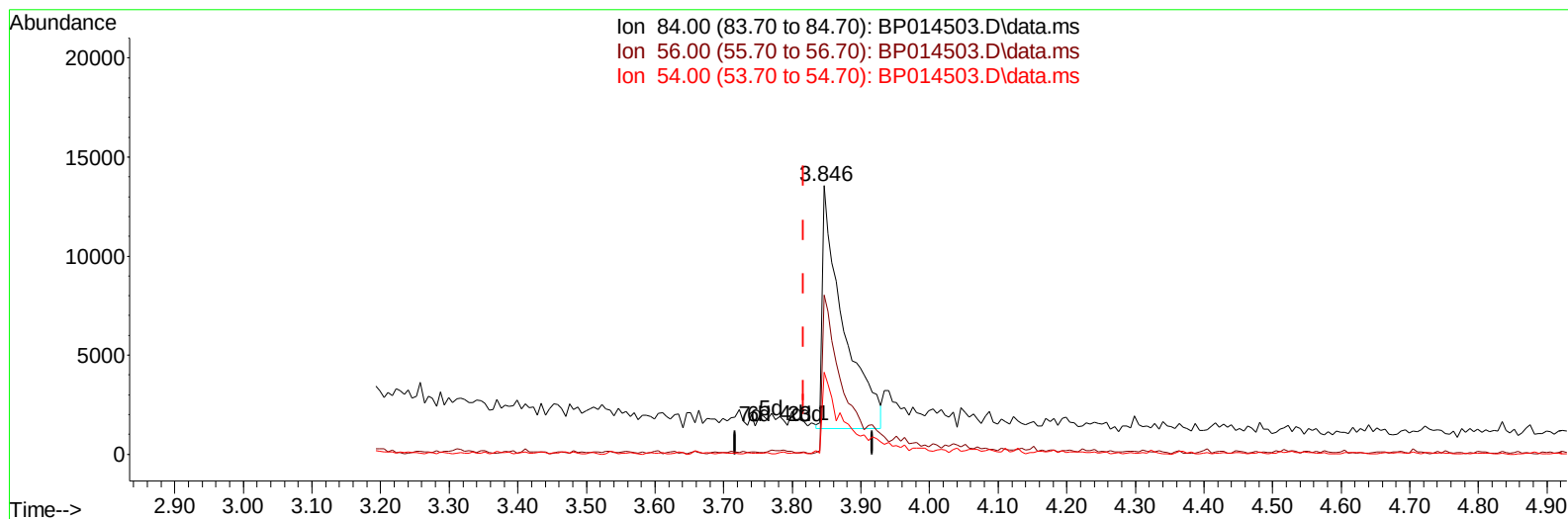
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(4) Pyridine-d5 (S)

3.846min (+ 0.029) 2.65 ng/ul m

response	25601	
Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	59.29
54.00	32.60	30.56
0.00	0.00	0.00

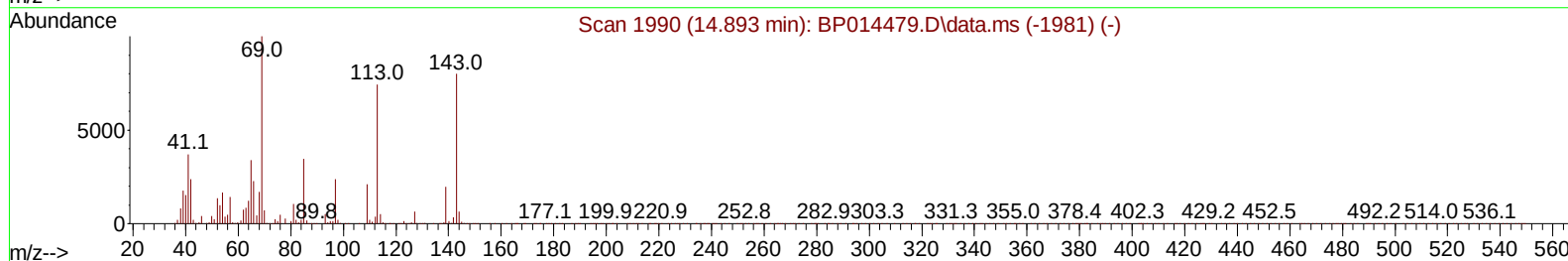
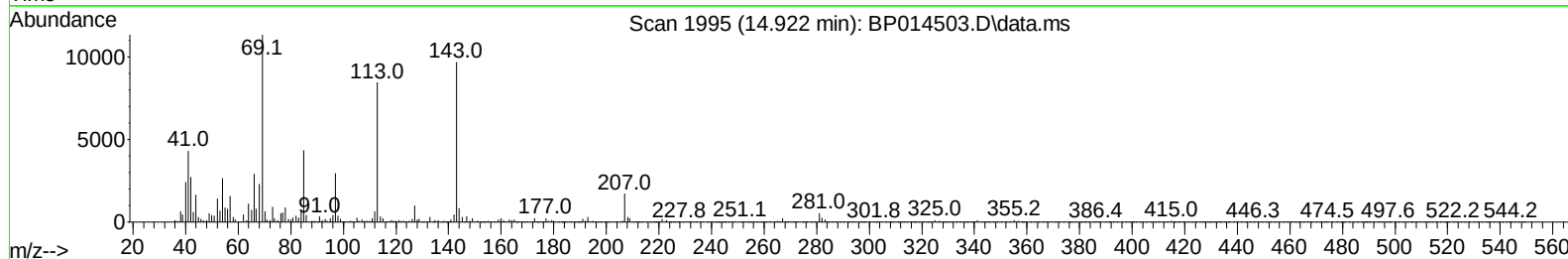
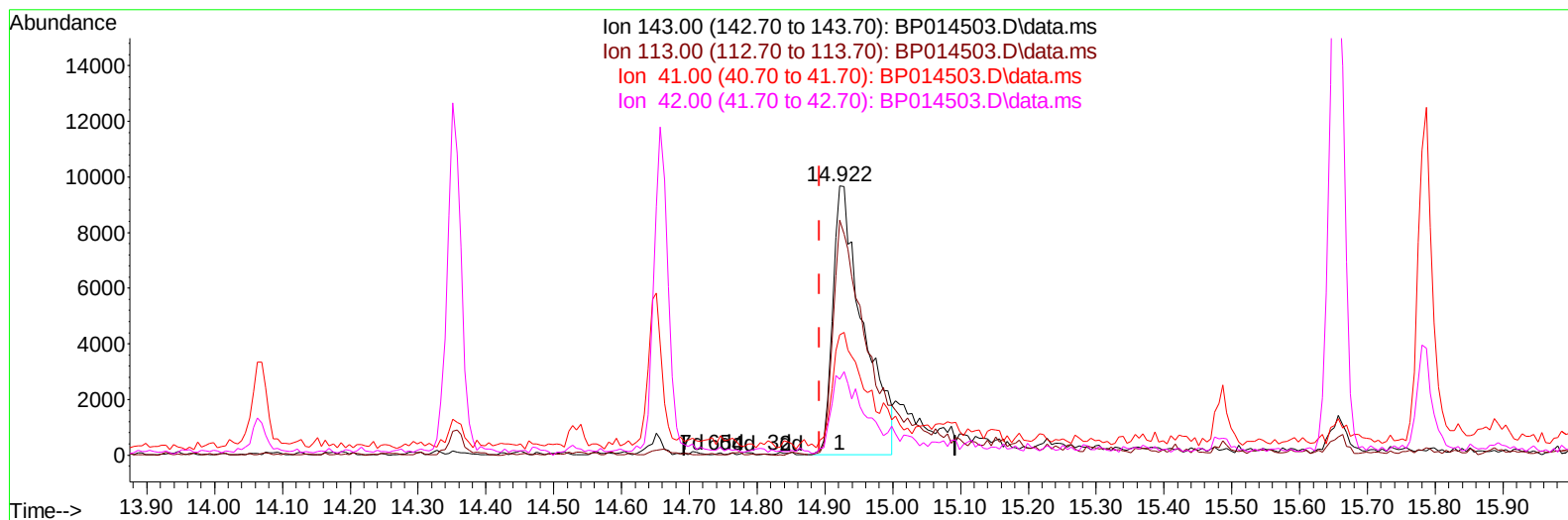
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TIC: BP014503.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.922min (+ 0.029) 5.30 ng/ul

response 29899

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	87.39
41.00	44.20	44.70
42.00	31.50	28.12

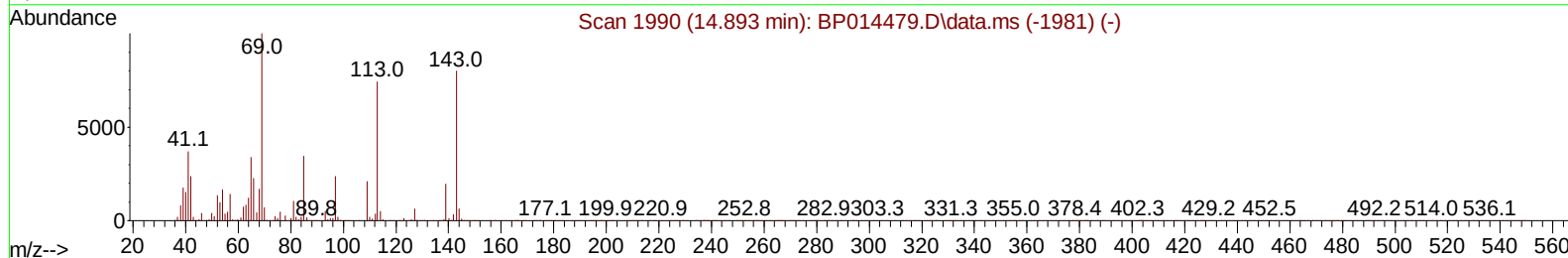
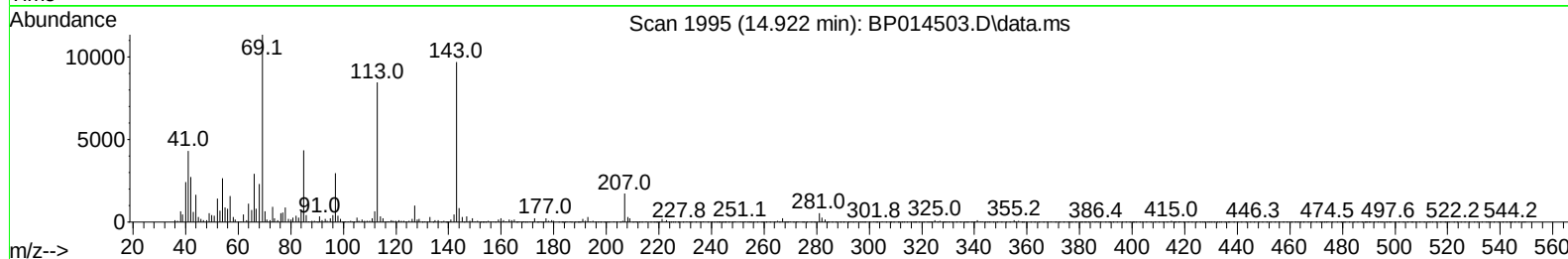
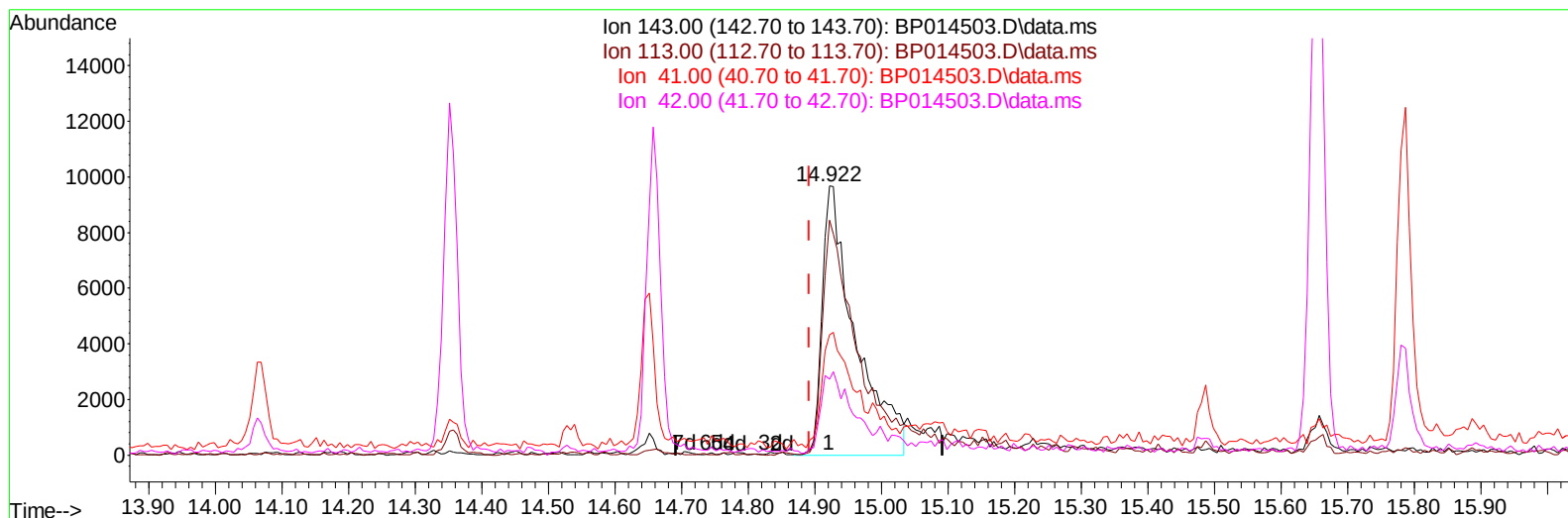
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Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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TIC: BP014503.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.922min (+ 0.029) 5.92 ng/ul m

response 33412

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	87.39
41.00	44.20	44.70
42.00	31.50	28.12

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.005	152	146185	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	658651	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.657	164	460899	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.422	188	1059223	20.000	ng/ul	0.00
79) Chrysene-d12	21.510	240	958544	20.000	ng/ul	0.00
88) Perylene-d12	24.021	264	913744	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	16152	4.285	ng/uL	0.00
4) Pyridine-d5	3.846	84	25601m	2.650	ng/ul	0.03
7) Phenol-d5	7.181	99	86186	6.393	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	254171	29.384	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	228671	23.467	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	168359	15.481	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	147193	27.683	ng/ul	0.00
24) 2-Nitrophenol-d4	9.910	143	170646	28.027	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.457	165	291726	26.309	ng/ul	0.00
31) 4-Chloroaniline-d4	10.981	131	169085	11.577	ng/ul	0.00
46) Dimethylphthalate-d6	14.069	166	1228219	32.866	ng/ul	0.00
49) Acenaphthylene-d8	14.357	160	1230038	29.663	ng/ul	0.00
54) 4-Nitrophenol-d4	14.922	143	33412m	5.918	ng/ul	0.03
60) Fluorene-d10	15.657	176	1049384	33.673	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.787	200	196468	25.208	ng/ul	0.00
73) Anthracene-d10	17.522	188	1867675	36.434	ng/ul	0.00
81) Pyrene-d10	19.751	212	2235232	34.717	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.863	264	1794366	36.962	ng/ul	0.00

Target Compounds Qvalue

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

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