

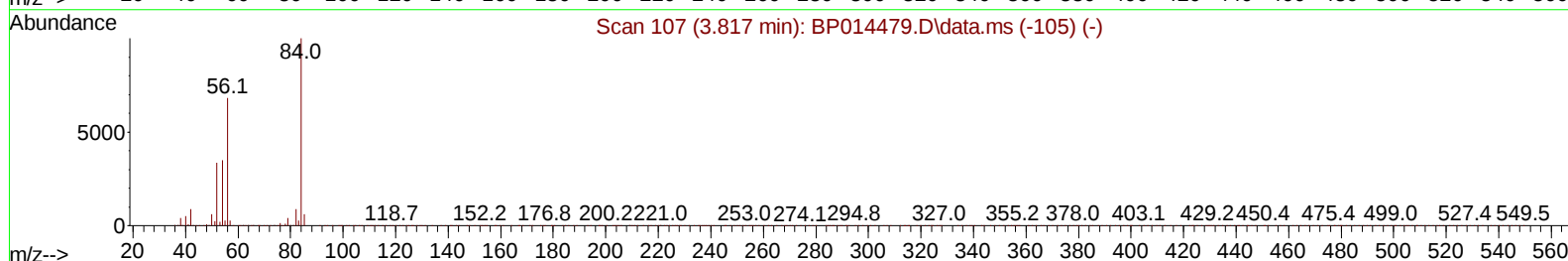
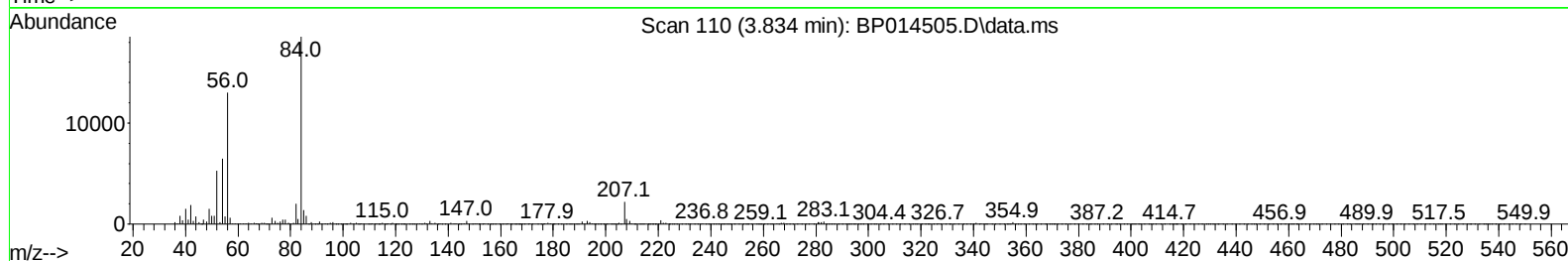
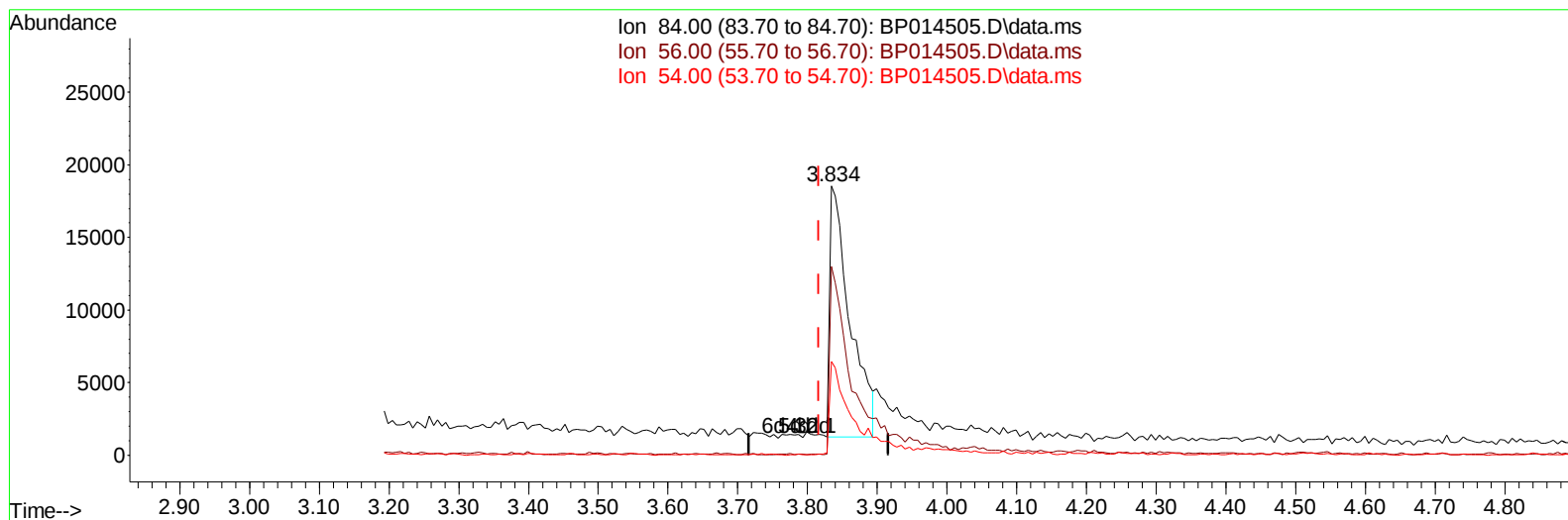
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
 Data File : BP014505.D
 Acq On : 04 Apr 2023 01:25
 Operator : CG/JU
 Sample : 02158-10
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PMW-1(20230328)

Manual IntegrationsAPPROVED

Quant Time: Apr 04 02:36:09 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: BP014505.D\data.ms

(4) Pyridine-d5 (S)

3.834min (+ 0.018) 4.22 ng/u1

response 34579

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	70.18
54.00	32.60	34.90
0.00	0.00	0.00

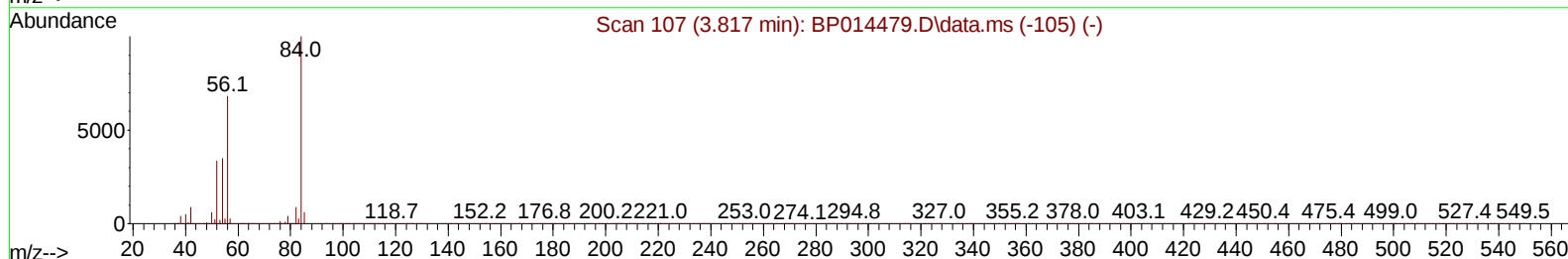
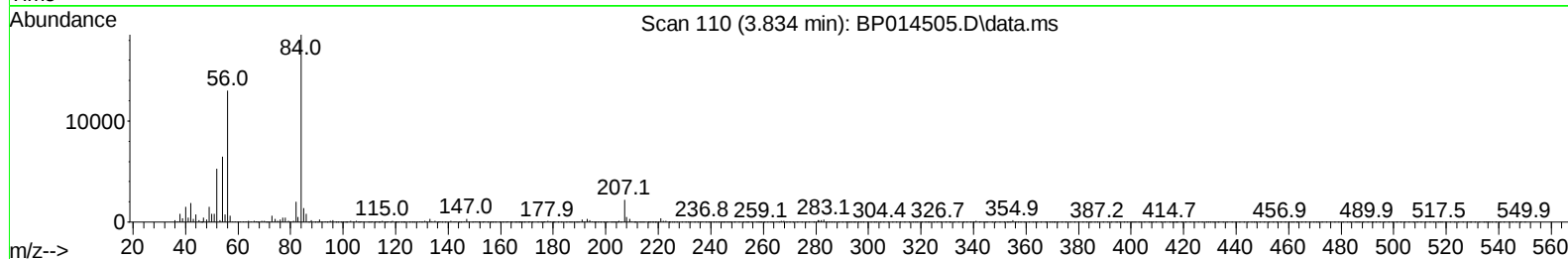
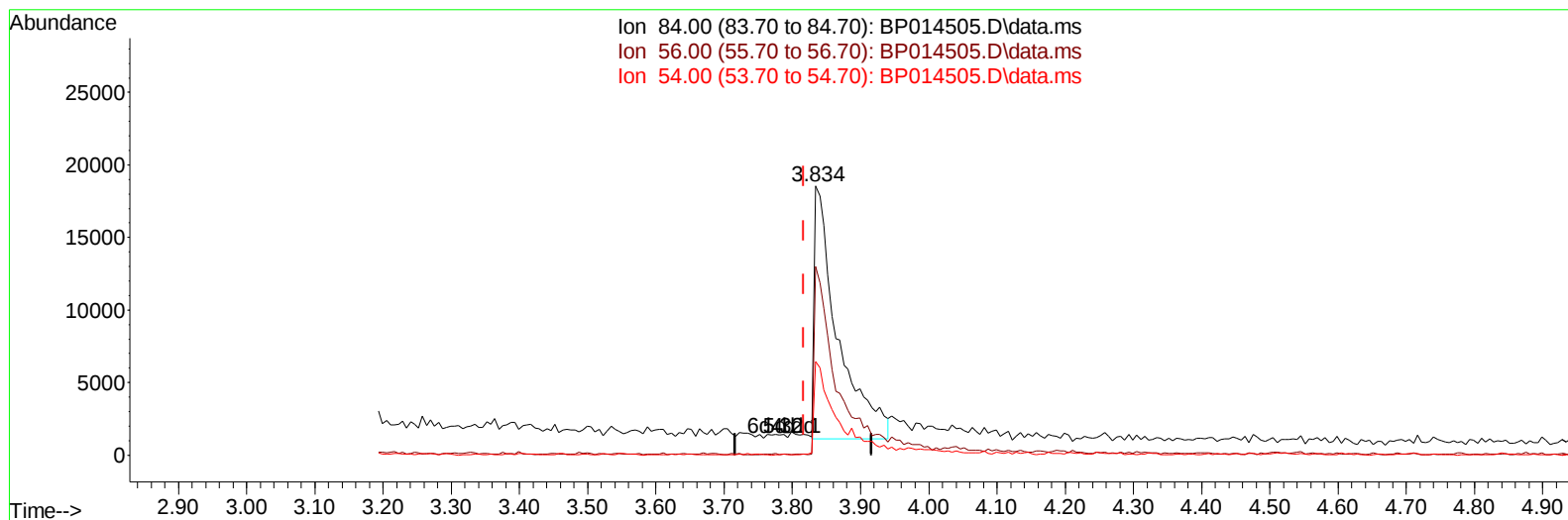
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(4) Pyridine-d5 (S)

3.834min (+ 0.018) 5.05 ng/ul m

response 41356

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	70.18
54.00	32.60	34.90
0.00	0.00	0.00

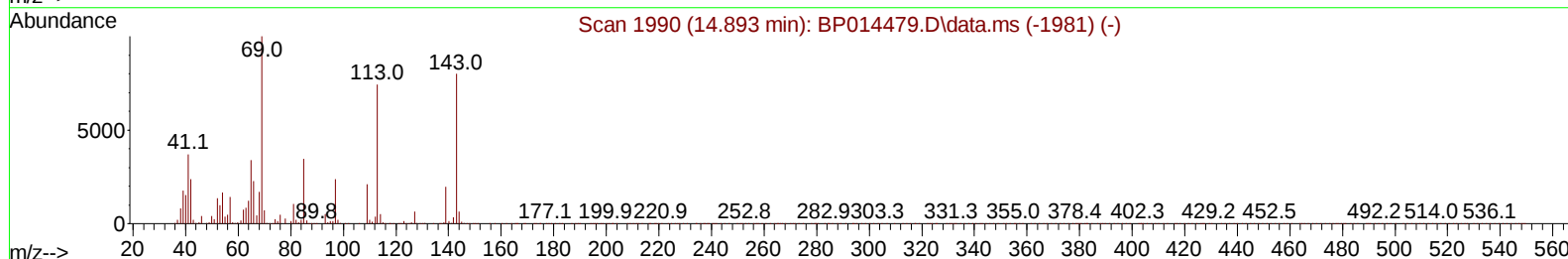
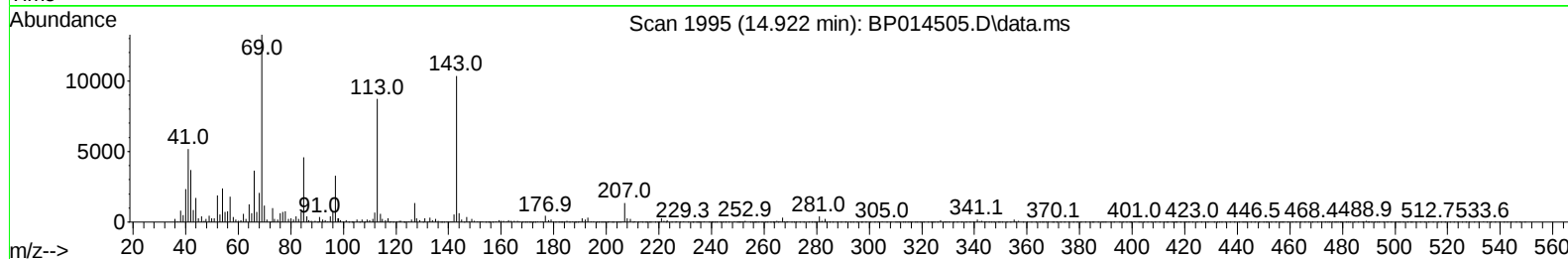
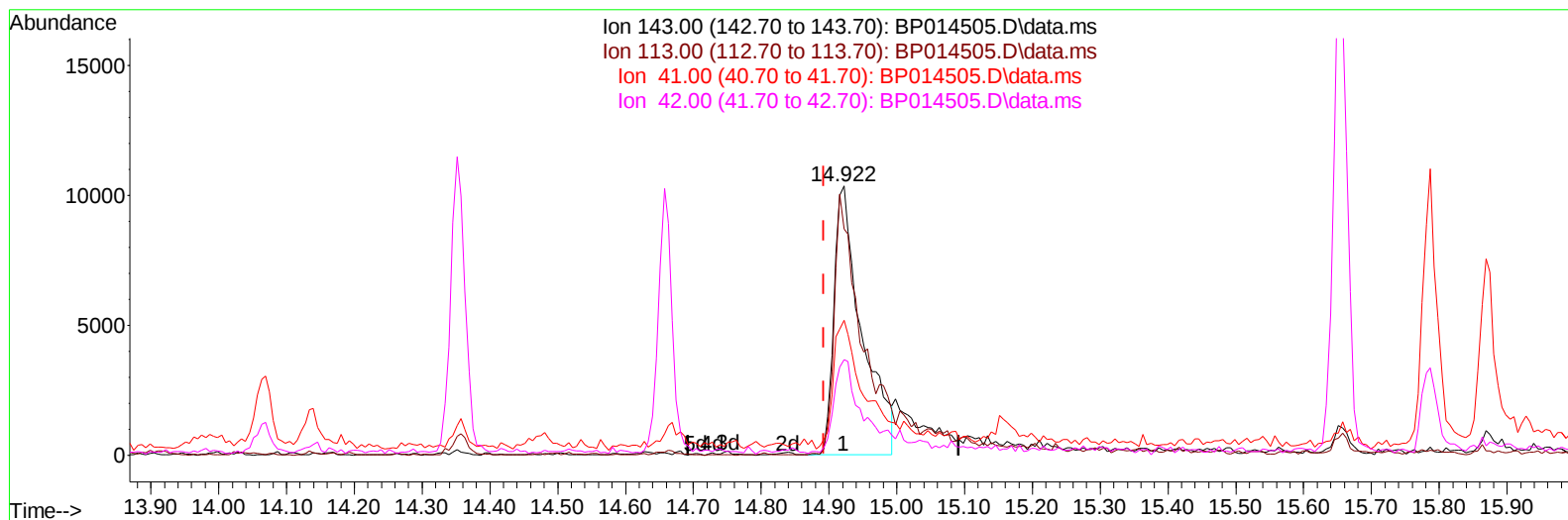
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TIC: BP014505.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.922min (+ 0.029) 6.21 ng/u1****response 29517**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	84.14
41.00	44.20	50.20
42.00	31.50	35.55

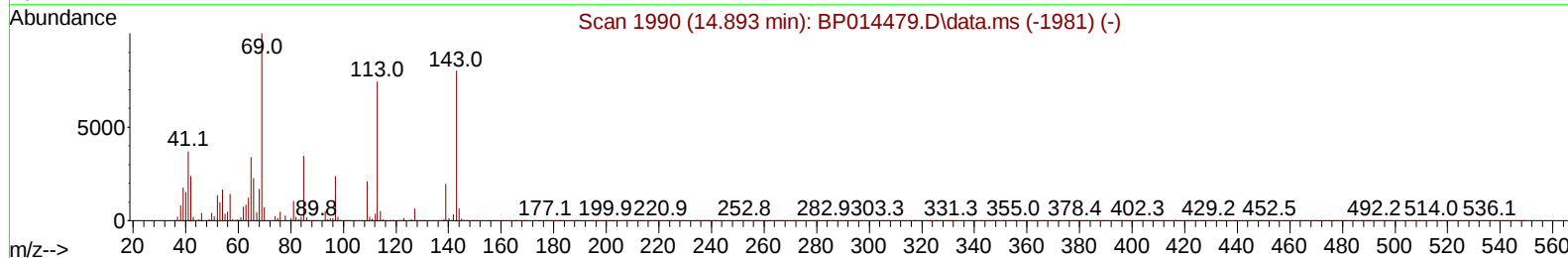
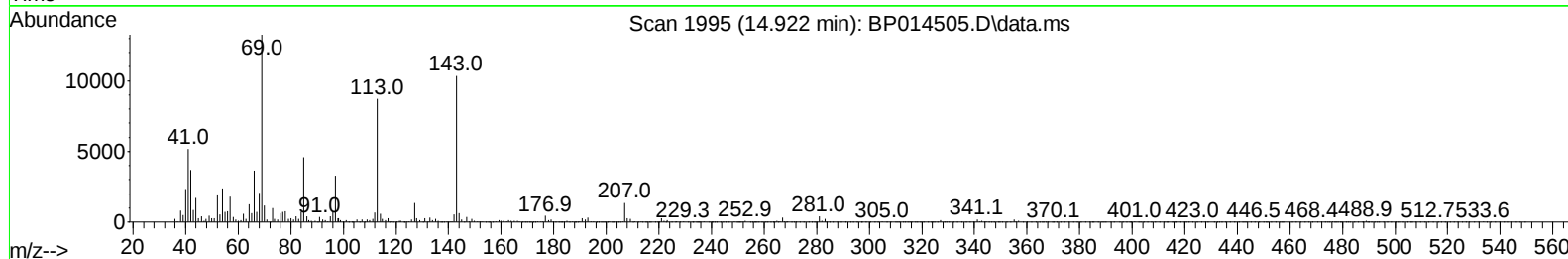
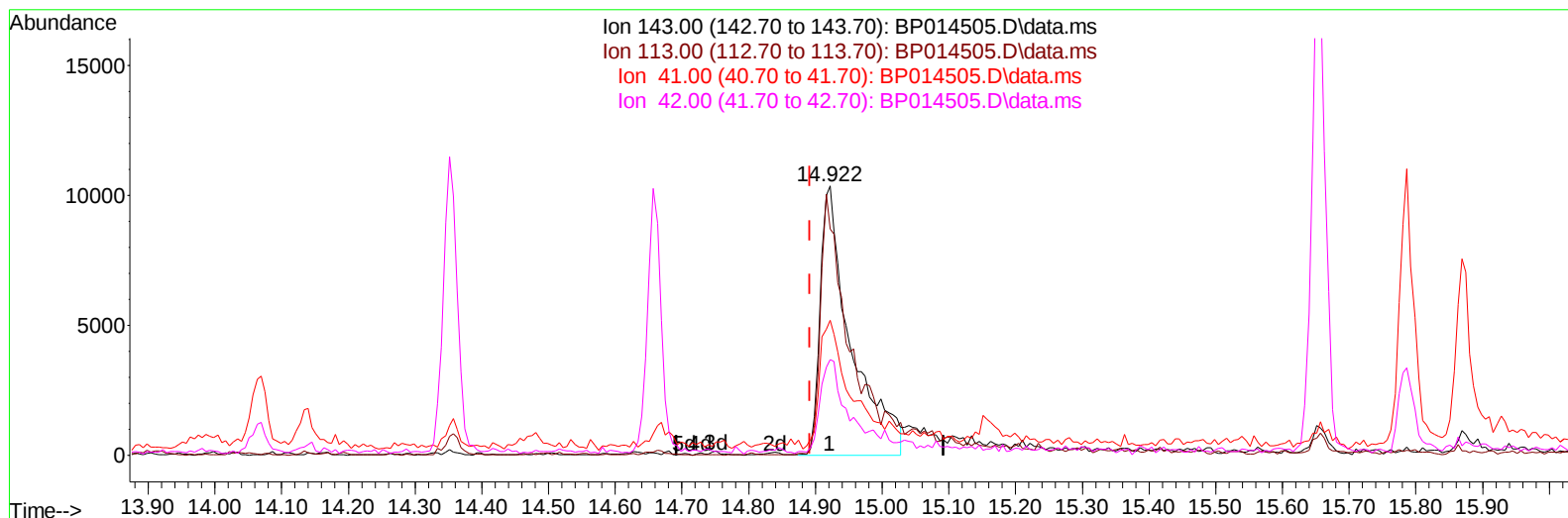
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 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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TIC: BP014505.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.922min (+ 0.029) 6.94 ng/ul m

response 32980

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	84.14
41.00	44.20	50.20
42.00	31.50	35.55

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.005	152	123885	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	565264	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.657	164	387735	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.422	188	932845	20.000	ng/ul	0.00
79) Chrysene-d12	21.510	240	935050	20.000	ng/ul	0.00
88) Perylene-d12	24.027	264	928998	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	14207	4.447	ng/uL	0.00
4) Pyridine-d5	3.834	84	41356m	5.052	ng/ul	0.02
7) Phenol-d5	7.181	99	93809	8.211	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	247303	33.736	ng/ul	0.00
11) 2-Chlorophenol-d4	7.534	132	229248	27.761	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	176119	19.110	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	149726	32.811	ng/ul	0.00
24) 2-Nitrophenol-d4	9.910	143	170753	32.678	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.457	165	274639	28.860	ng/ul	0.00
31) 4-Chloroaniline-d4	10.981	131	206900	16.506	ng/ul	0.00
46) Dimethylphthalate-d6	14.069	166	1090070	34.673	ng/ul	0.00
49) Acenaphthylene-d8	14.351	160	1185916	33.996	ng/ul	0.00
54) 4-Nitrophenol-d4	14.922	143	32980m	6.944	ng/ul	0.03
60) Fluorene-d10	15.657	176	922970	35.205	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.787	200	172050	25.066	ng/ul	0.00
73) Anthracene-d10	17.522	188	1602345	35.493	ng/ul	0.00
81) Pyrene-d10	19.751	212	1963961	31.270	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.863	264	1759199	35.642	ng/ul	0.00

Target Compounds Qvalue

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

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