

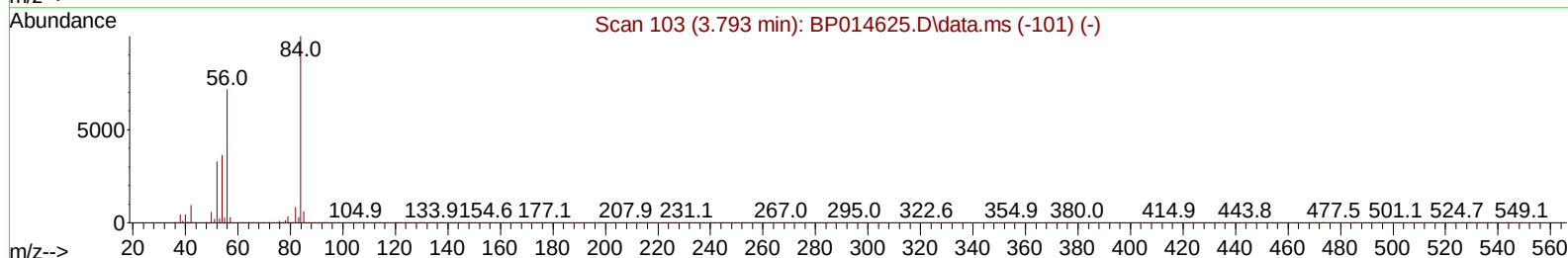
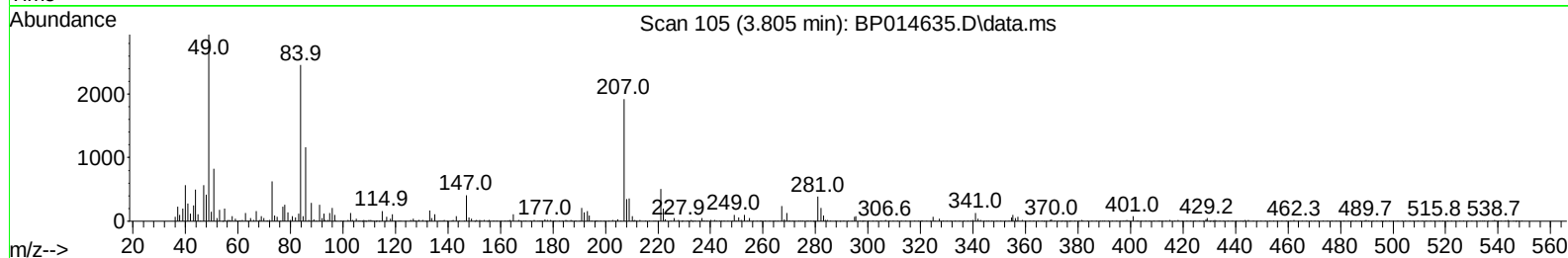
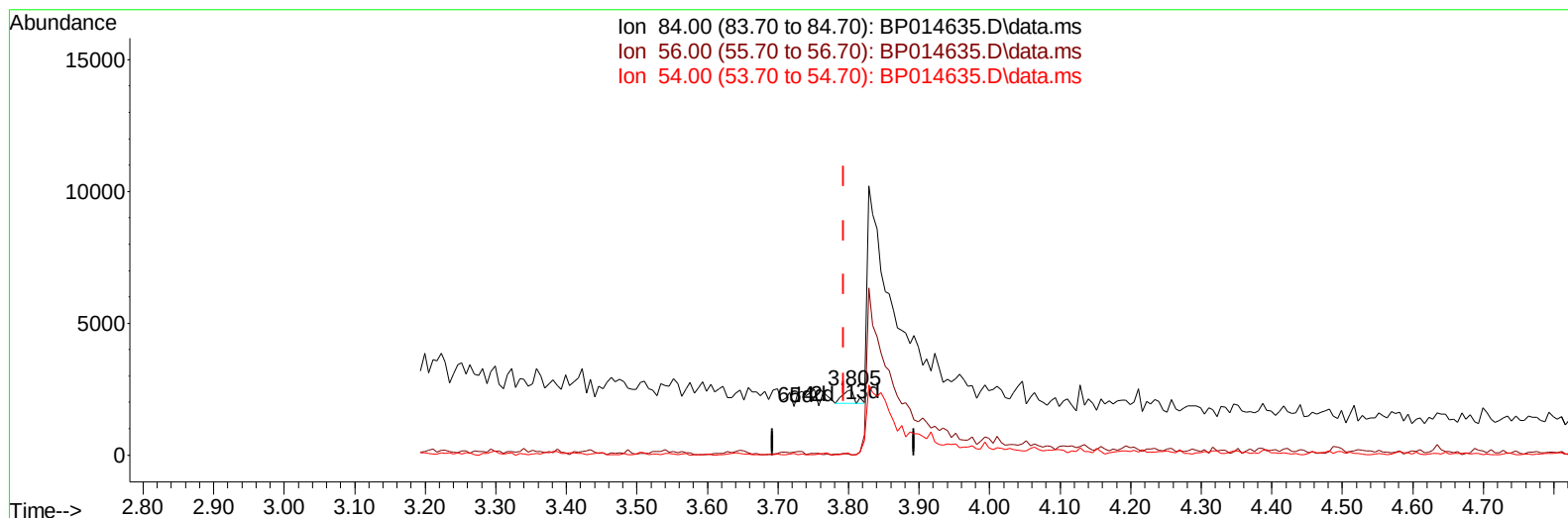
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041023\
Data File : BP014635.D
Acq On : 10 Apr 2023 19:56
Operator : CG/JU
Sample : 02250-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHDx8

Manual IntegrationsAPPROVED

Quant Time: Apr 11 01:24:18 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Apr 11 01:21:43 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/11/2023
Supervised By :Jagrut Upadhyay 04/11/2023



TIC: BP014635.D\data.ms

(4) Pyridine-d5 (S)

3.805min (+ 0.012) 0.09 ng/u1

response 633

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	0.86#
54.00	32.60	0.08#
0.00	0.00	0.00

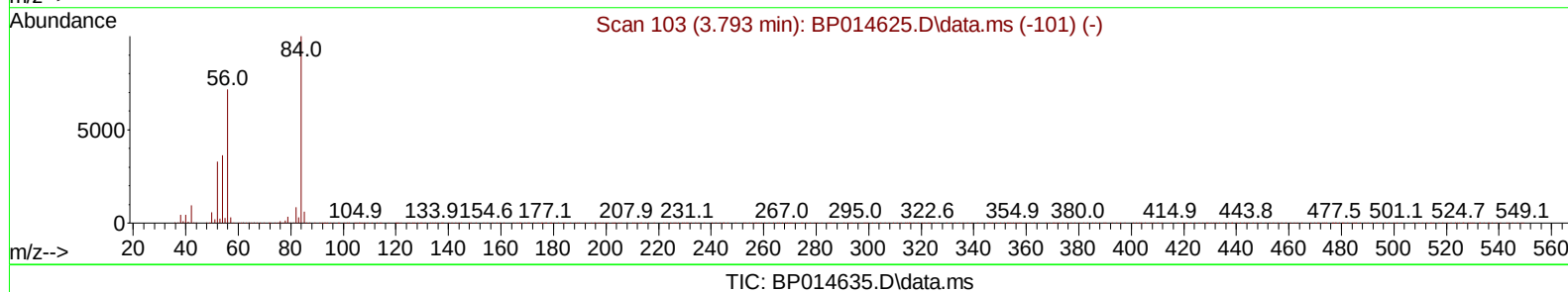
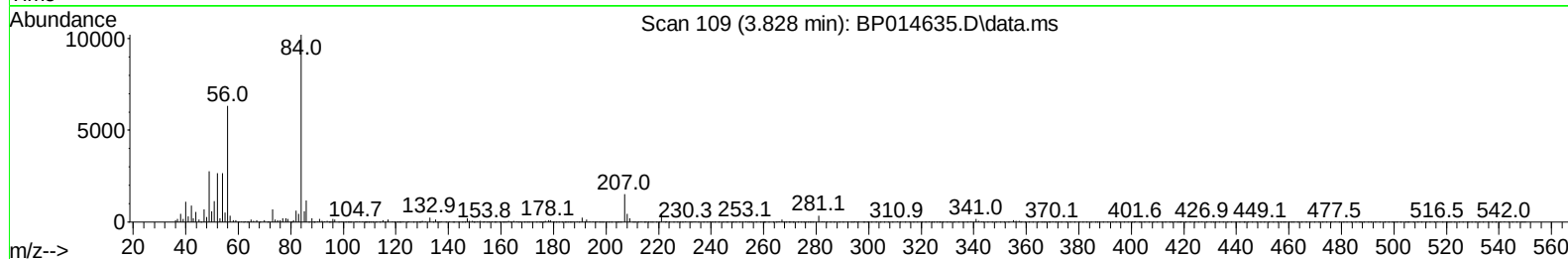
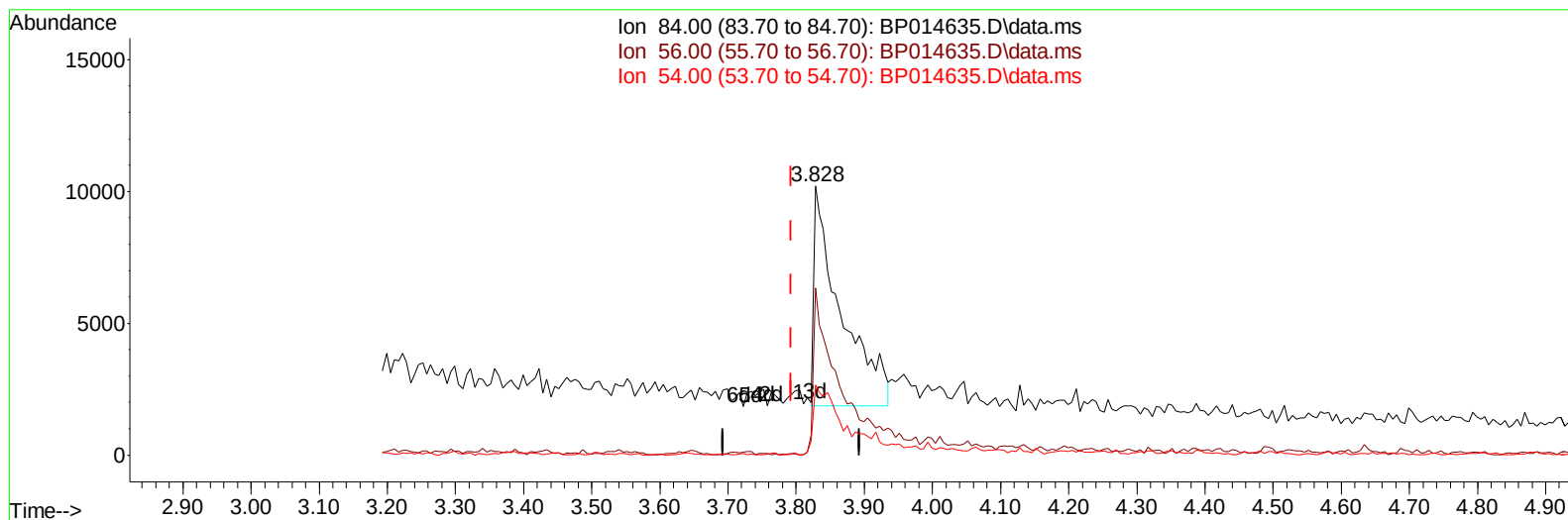
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(4) Pyridine-d5 (S)

3.828min (+ 0.035) 3.36 ng/ul m

response 22773

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	62.04
54.00	32.60	26.00#
0.00	0.00	0.00

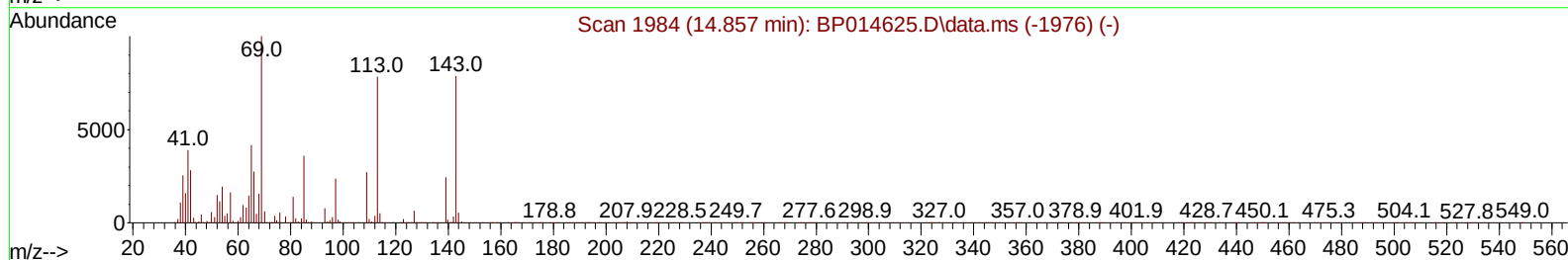
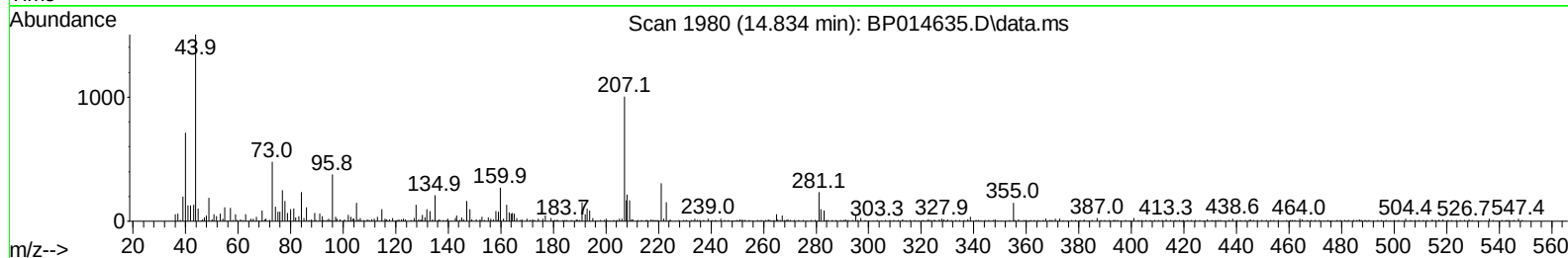
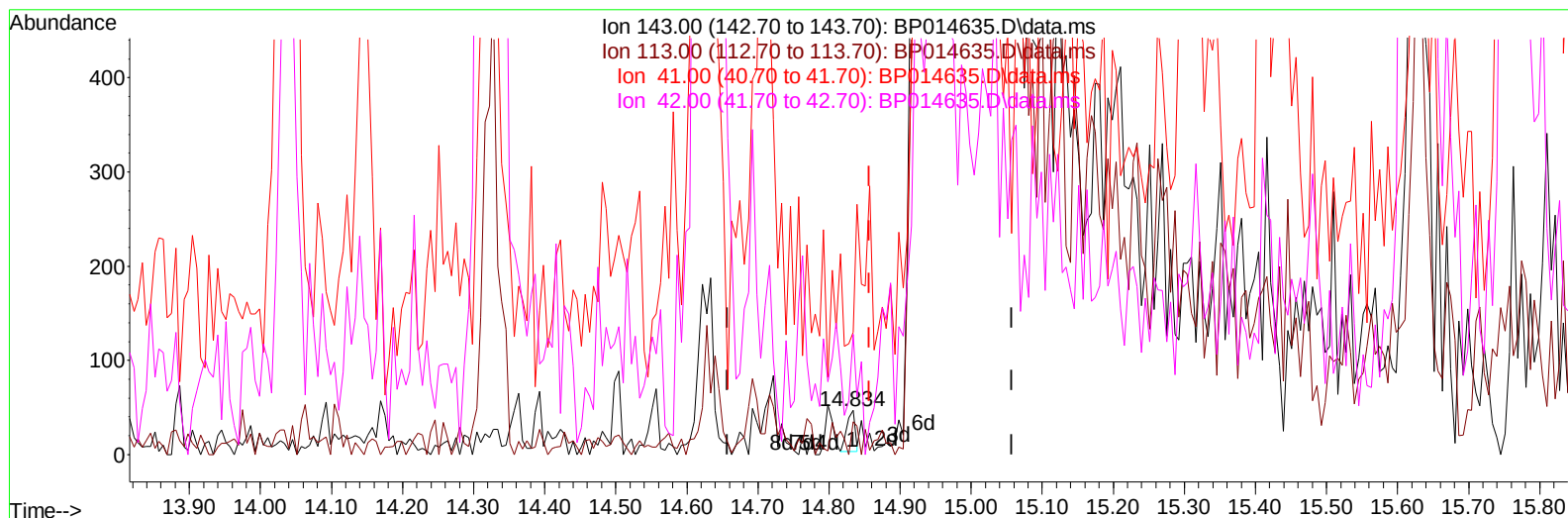
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TIC: BP014635.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.834min (-0.023) 0.01 ng/ul

response 35

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	74.47#
41.00	44.20	276.60#
42.00	31.50	272.34#

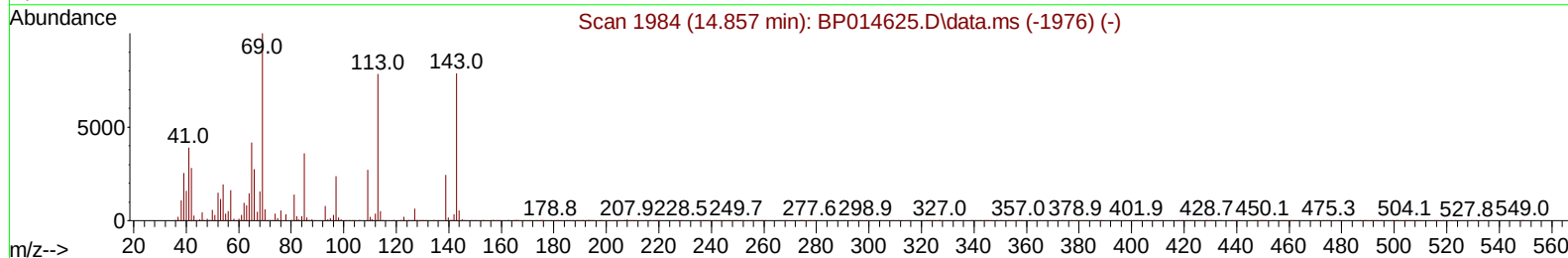
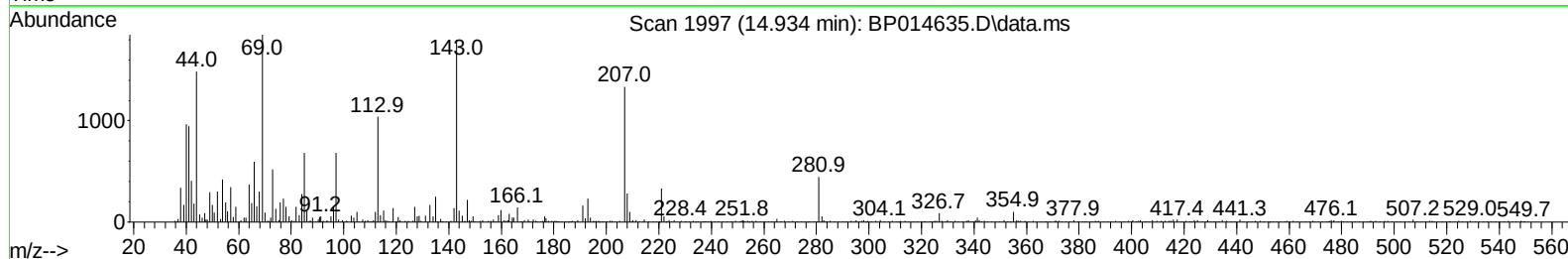
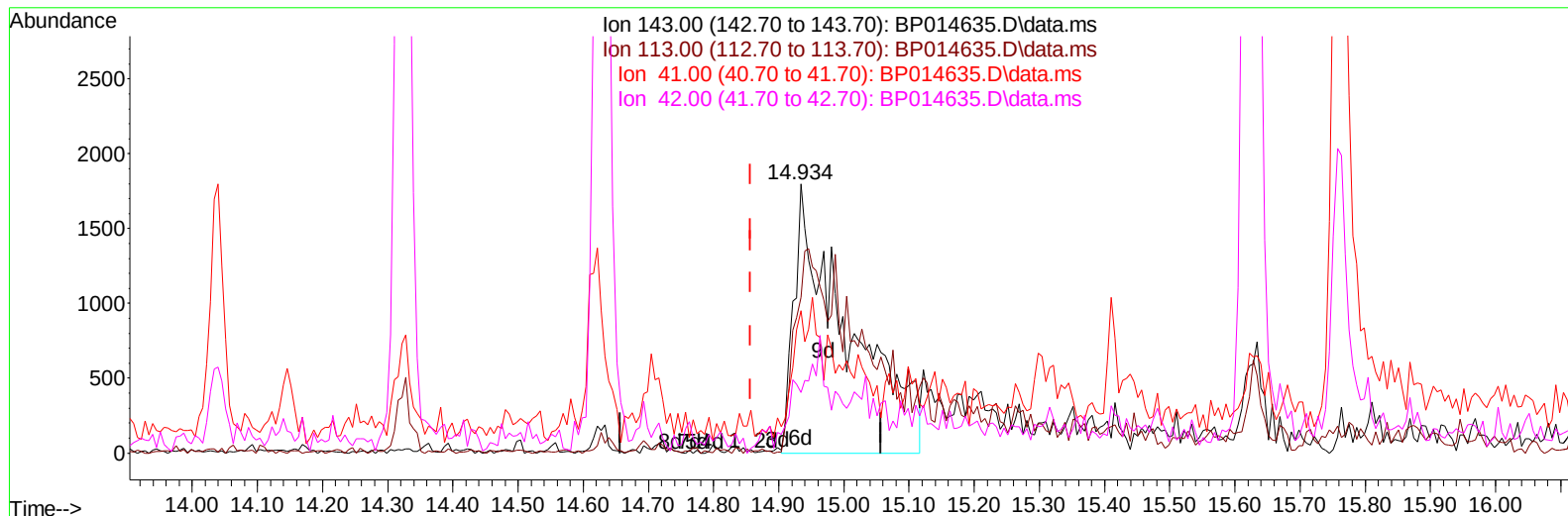
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TIC: BP014635.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.934min (+ 0.076) 3.23 ng/ul m

response 10216

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	57.87#
41.00	44.20	52.81
42.00	31.50	22.59#

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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	7.975	152	102569	20.000	ng/ul	0.00
20) Naphthalene-d8	10.799	136	424751	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.628	164	257867	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	553227	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	559524	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	644217	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	12305	4.652	ng/uL	0.00
4) Pyridine-d5	3.828	84	22773m	3.360	ng/ul	0.04
7) Phenol-d5	7.158	99	48947	5.174	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.305	67	162893	26.839	ng/ul	0.00
11) 2-Chlorophenol-d4	7.510	132	142775	20.883	ng/ul	0.00
15) 4-Methylphenol-d8	8.705	113	92153	12.077	ng/ul	0.00
21) Nitrobenzene-d5	9.157	128	82913	24.180	ng/ul	0.00
24) 2-Nitrophenol-d4	9.881	143	90653	23.088	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.428	165	143702	20.096	ng/ul	0.00
31) 4-Chloroaniline-d4	10.951	131	159211	16.903	ng/ul	0.00
46) Dimethylphthalate-d6	14.040	166	565152	27.030	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	605102	26.082	ng/ul	0.00
54) 4-Nitrophenol-d4	14.934	143	10216m	3.234	ng/ul	0.08
60) Fluorene-d10	15.628	176	480999	27.587	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	86367	21.217	ng/ul	0.00
73) Anthracene-d10	17.492	188	891940	33.314	ng/ul	0.00
81) Pyrene-d10	19.722	212	1084859	28.866	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	1177646	34.407	ng/ul	0.00

Target Compounds Qvalue

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

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