

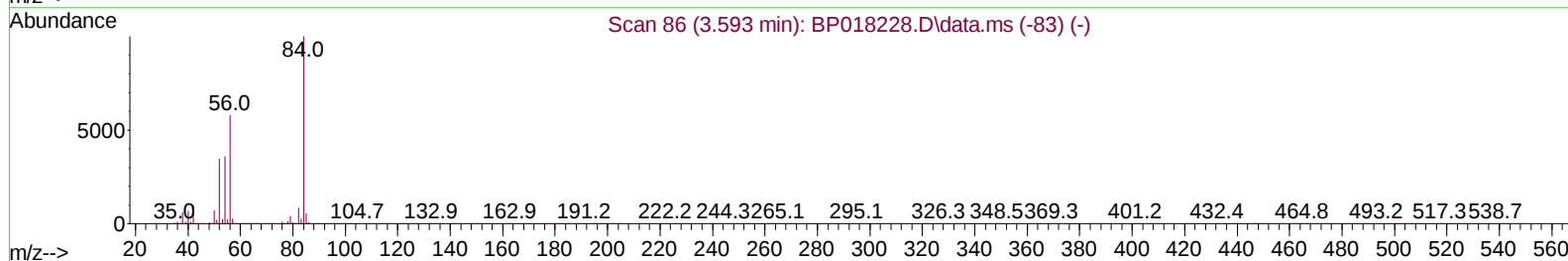
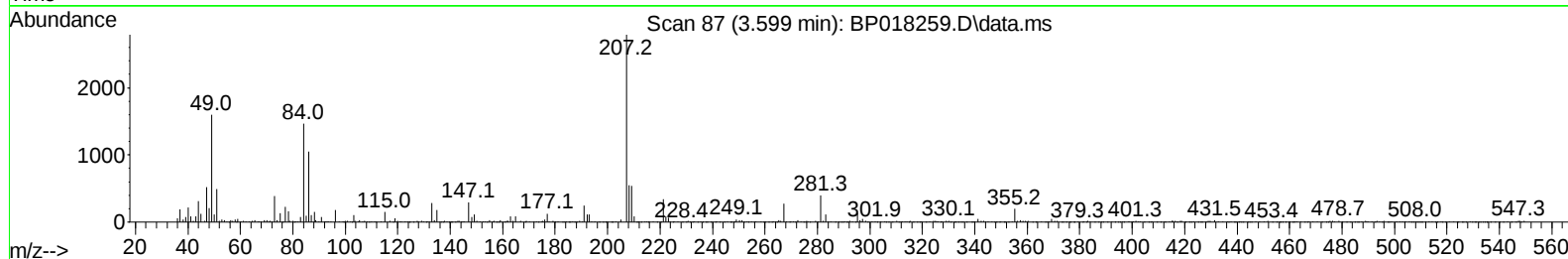
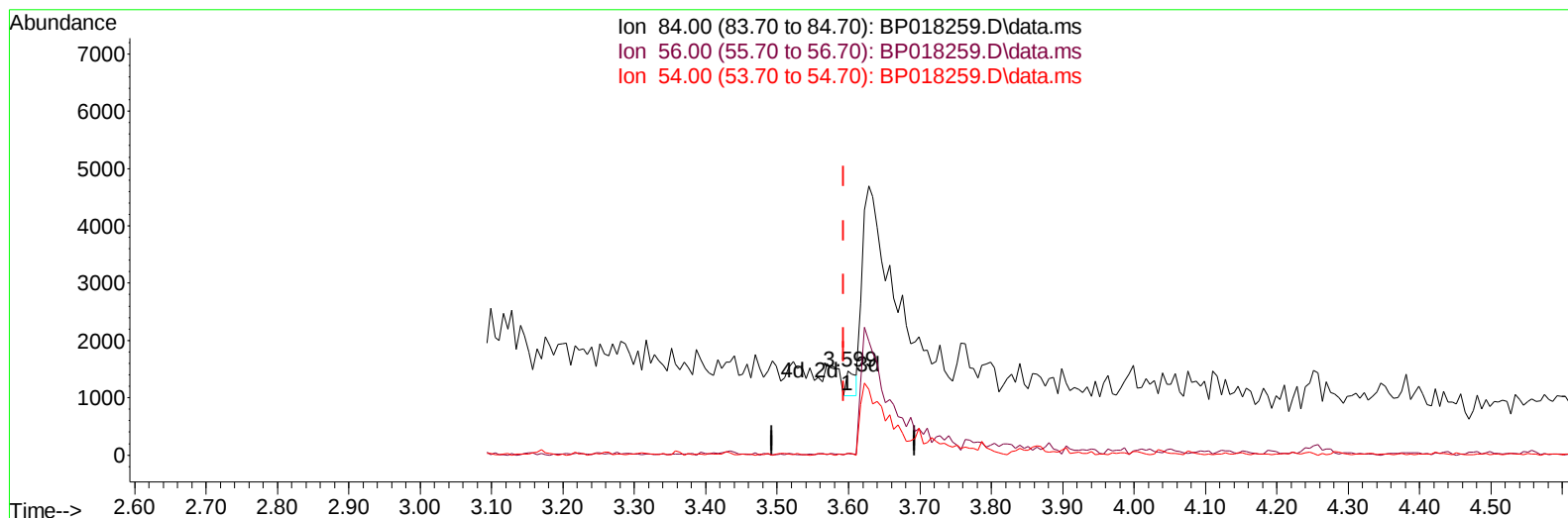
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
 Data File : BP018259.D
 Acq On : 18 Nov 2023 13:49
 Operator : MA/JU
 Sample : 05370-14
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDQ8

Manual IntegrationsAPPROVED

Quant Time: Nov 19 00:52:50 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 18:13:37 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/19/2023
 Supervised By :mohammad ahmed 11/19/2023



TIC: BP018259.D\data.ms

(4) Pyridine-d5 (S)

3.599min (+ 0.006) 0.14 ng/u1

response 401

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.30	1.71#
54.00	35.20	1.71#
0.00	0.00	0.00

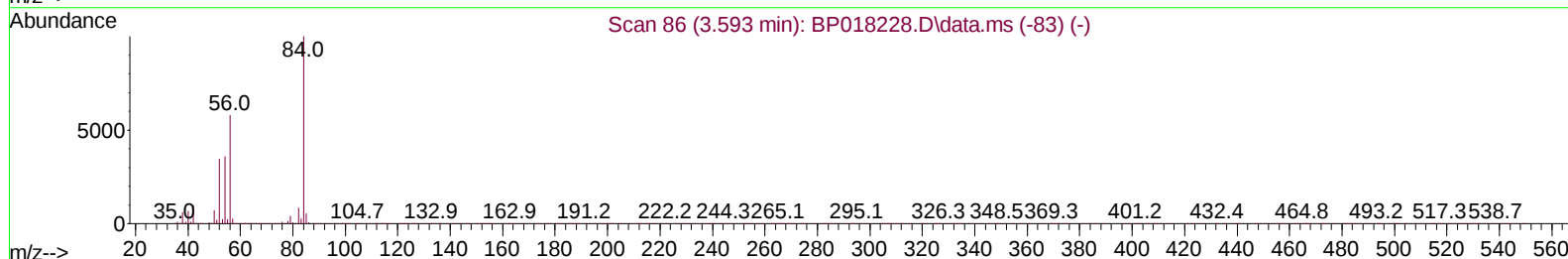
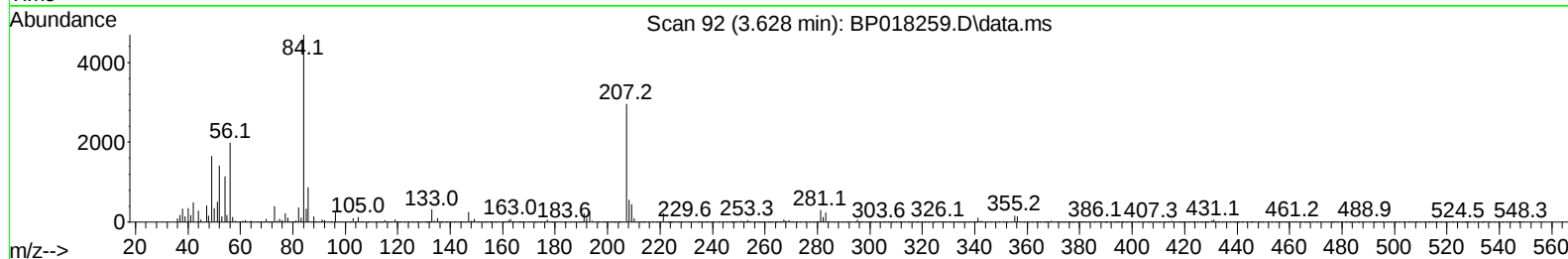
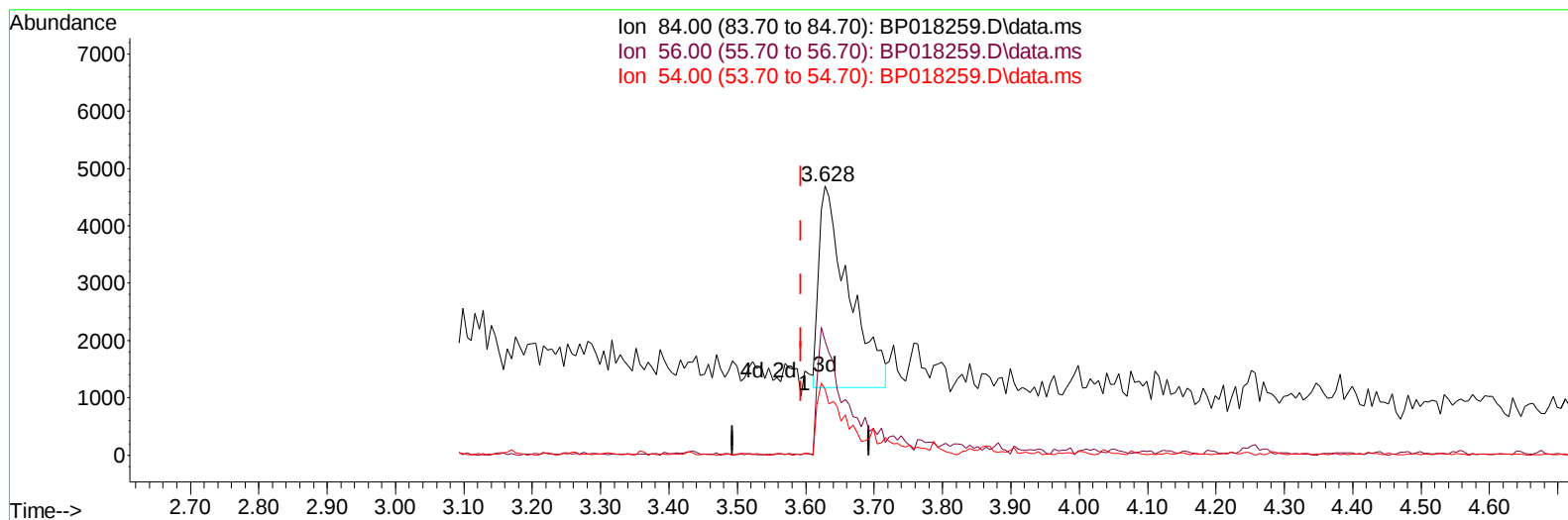
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(4) Pyridine-d5 (S)

3.628min (+ 0.035) 3.77 ng/ul m

response 10624

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.30	42.30#
54.00	35.20	24.53#
0.00	0.00	0.00

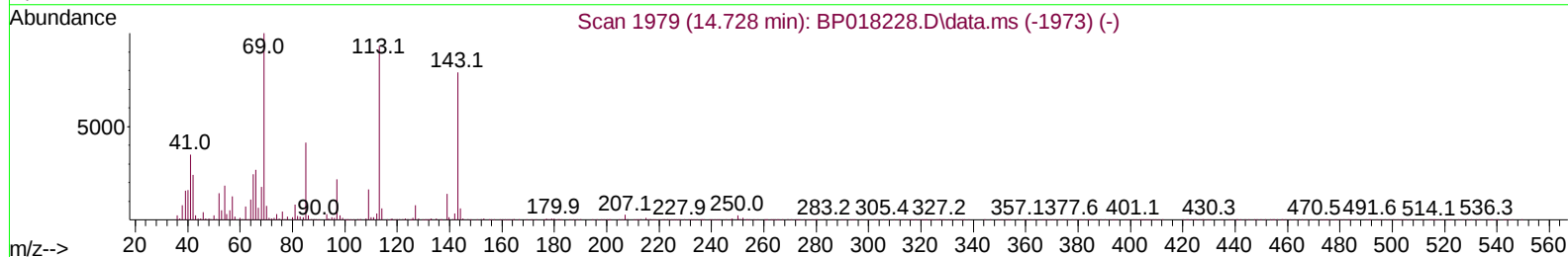
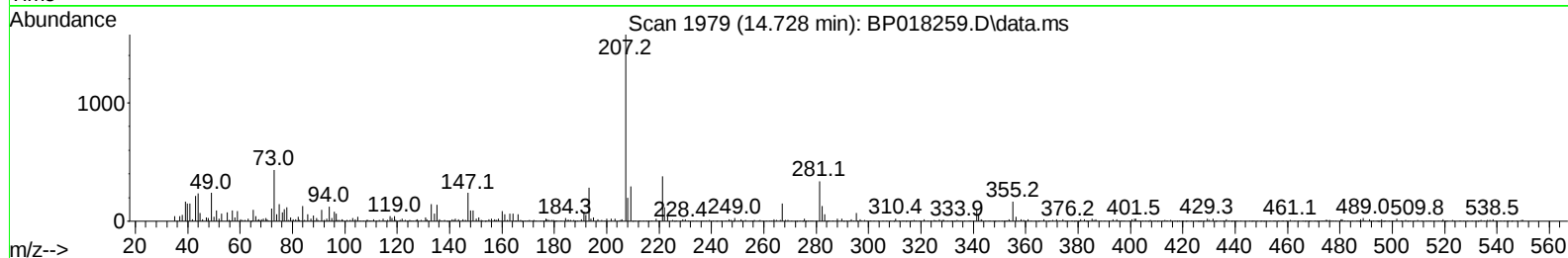
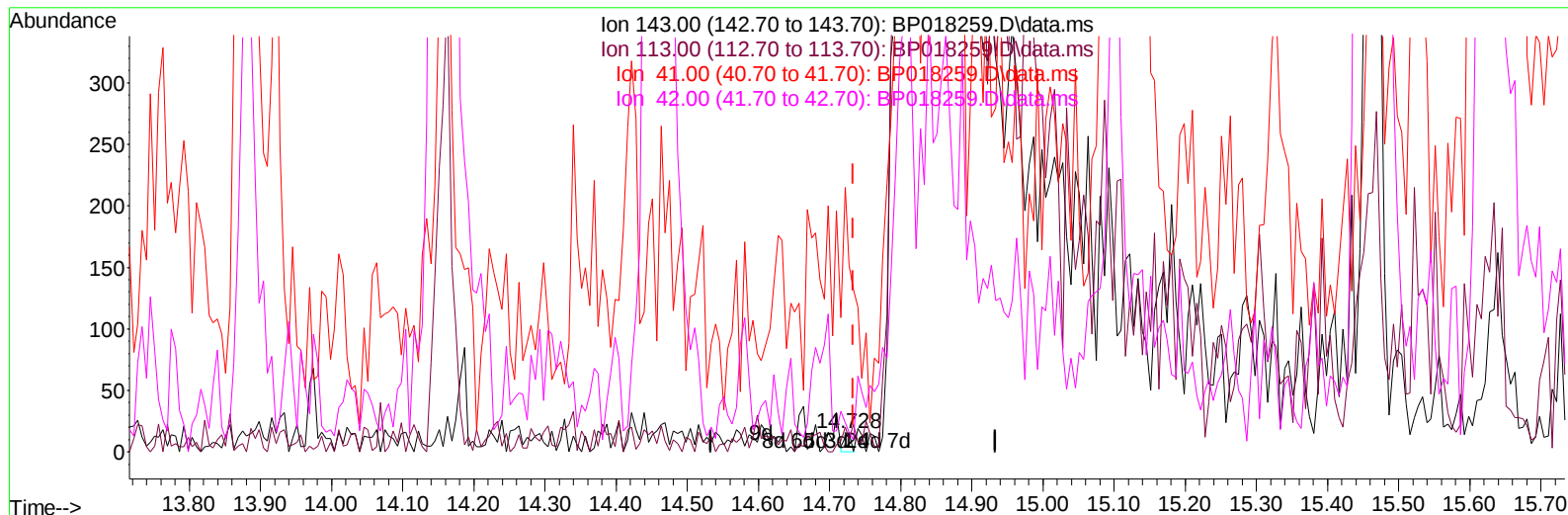
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TIC: BP018259.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.728min (-0.006) 0.01 ng/u1

response 10

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	68.75#
41.00	47.70	950.00#
42.00	30.10	100.00#

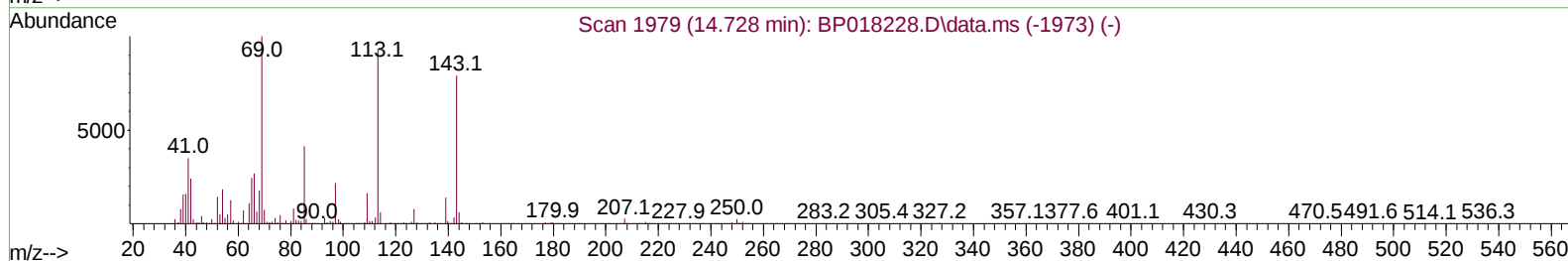
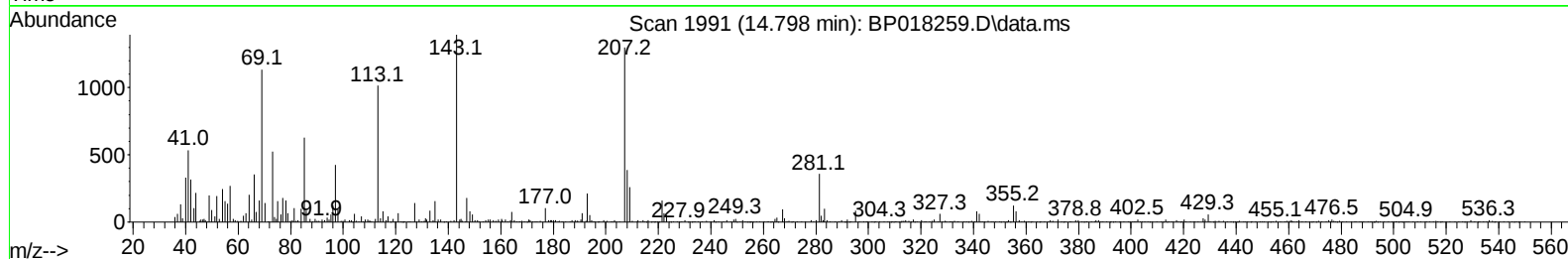
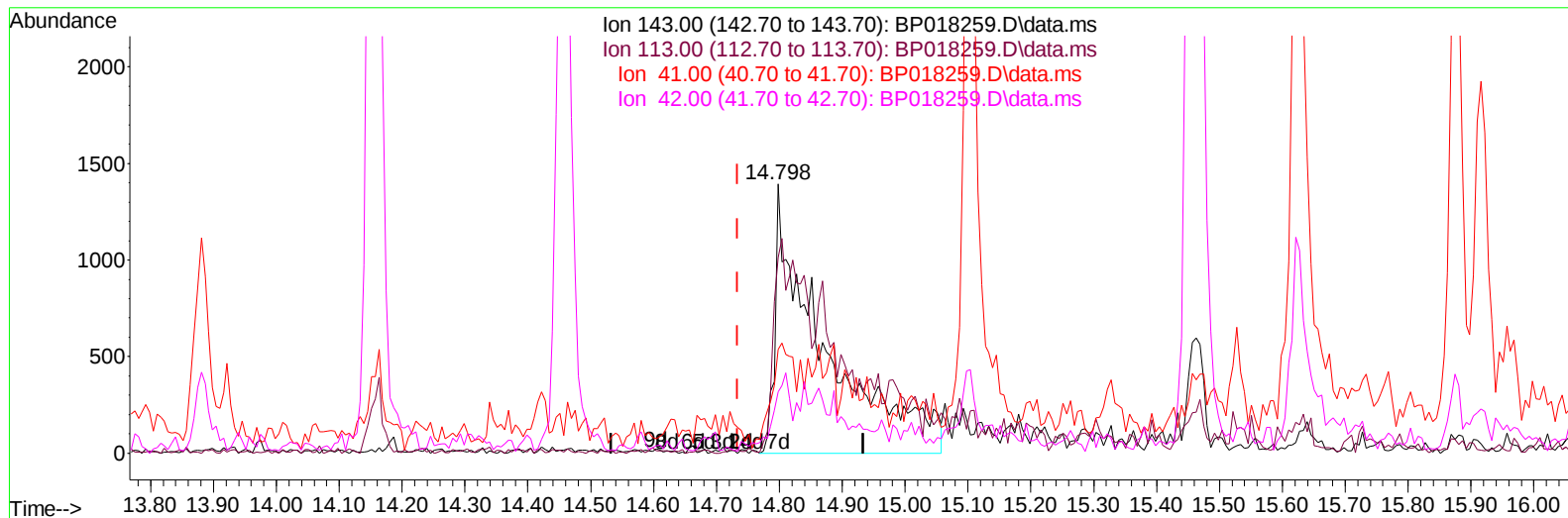
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TIC: BP018259.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.798min (+ 0.065) 5.46 ng/ul m

response 7353

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	72.79#
41.00	47.70	38.48
42.00	30.10	22.83#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	38110	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	157037	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	97753	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.239	188	250573	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	256762	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	286922	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	5403	5.001	ng/uL	0.00
4) Pyridine-d5	3.628	84	10624m	3.772	ng/ul	0.04
7) Phenol-d5	6.963	99	21263	5.779	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.128	67	74771	34.581	ng/ul	0.00
11) 2-Chlorophenol-d4	7.305	132	74987	25.737	ng/ul	0.00
15) 4-Methylphenol-d8	8.510	113	50789	16.846	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	48179	34.053	ng/ul	0.00
24) 2-Nitrophenol-d4	9.698	143	51311	32.012	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	85610	29.967	ng/ul	0.00
31) 4-Chloroaniline-d4	10.781	131	109578	27.970	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	392535	43.521	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	374680	38.140	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	7353m	5.456	ng/ul	0.06
60) Fluorene-d10	15.463	176	319336	42.884	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	53794	29.697	ng/ul	0.00
73) Anthracene-d10	17.339	188	583202	43.034	ng/ul	0.00
81) Pyrene-d10	19.598	212	754824	44.011	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.686	264	760037	45.556	ng/ul	0.00

Target Compounds Qvalue

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

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