

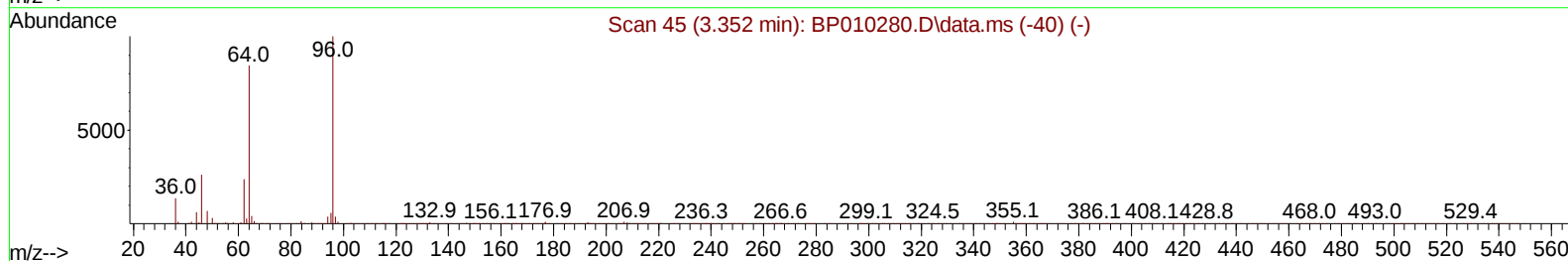
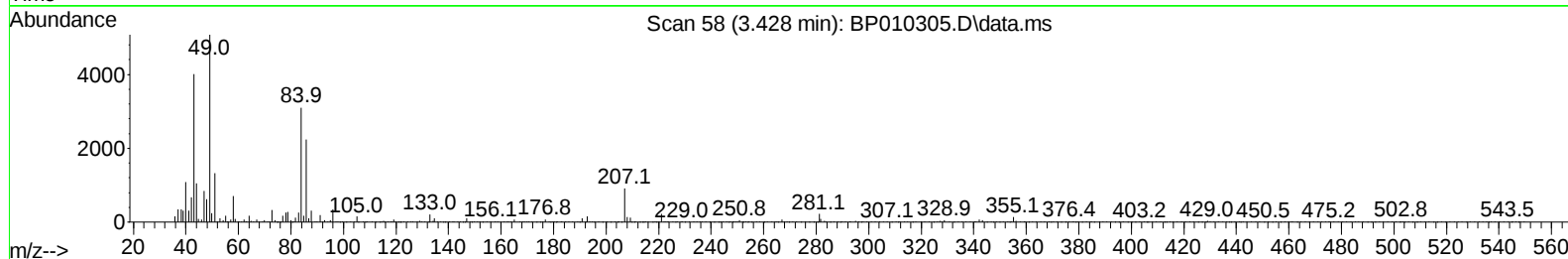
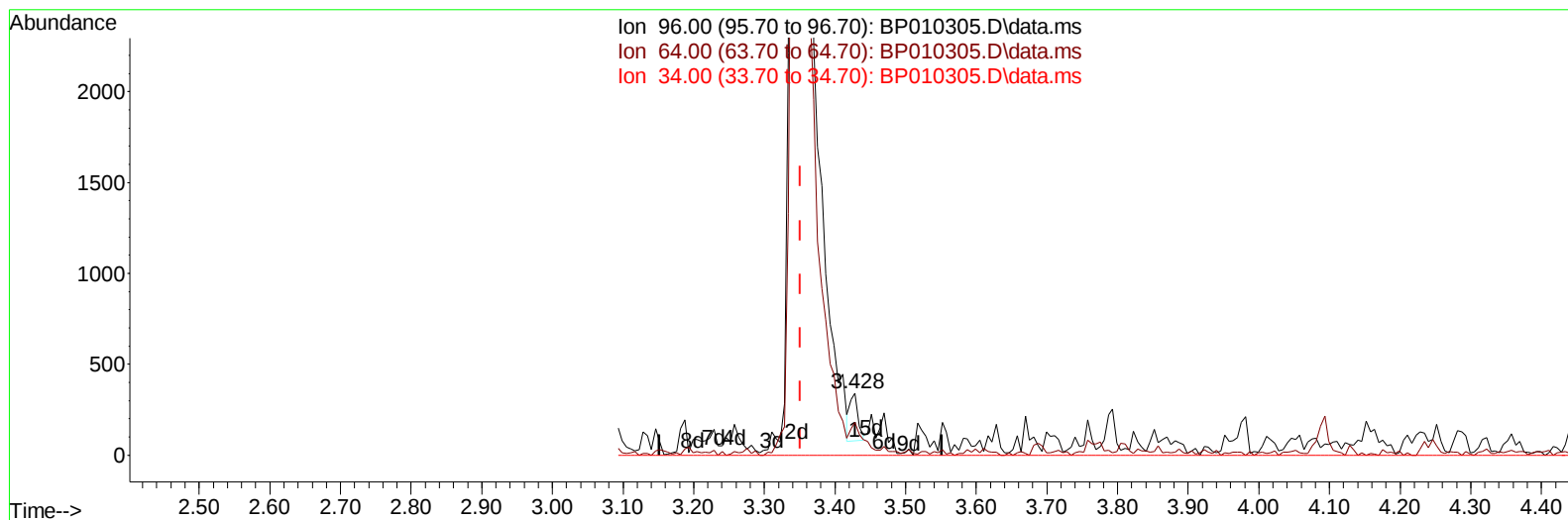
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010305.D
Acq On : 12 May 2022 02:43
Operator : CG/JU
Sample : N2707-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXL85

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/12/2022
Supervised By :Yogesh Patel 05/17/2022

Quant Time: May 12 05:33:06 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu May 12 01:10:15 2022
Response via : Initial Calibration



TIC: BP010305.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.428min (+ 0.077) 0.07 ng/uL

response 211

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	64.52
34.00	0.00	0.00
0.00	0.00	0.00

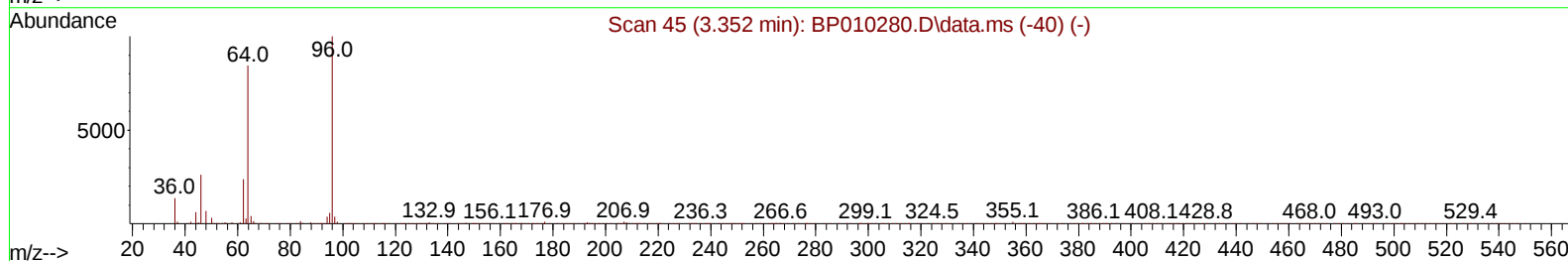
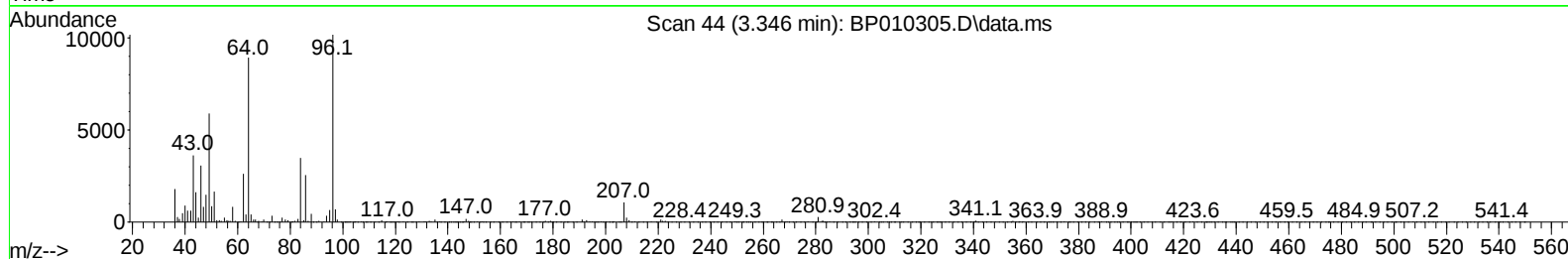
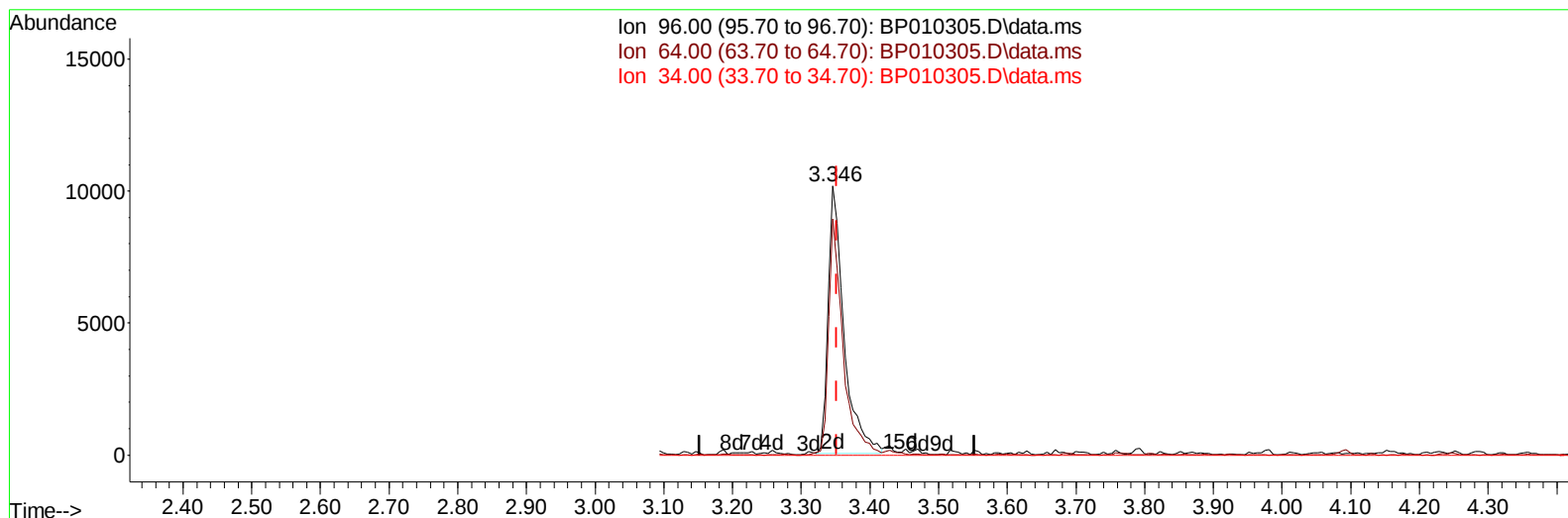
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(3) 1,4-Dioxane-d8 (S)

3.346min (-0.006) 5.01 ng/uL m

response 16235

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	87.72#
34.00	0.00	0.00
0.00	0.00	0.00

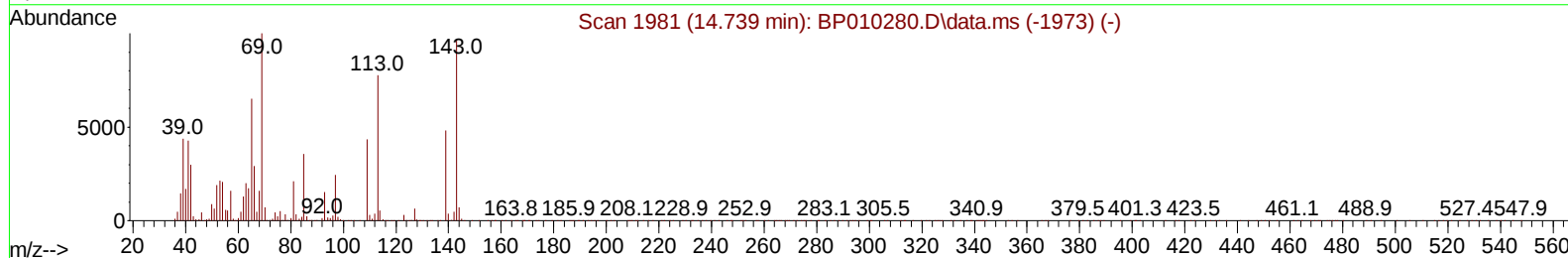
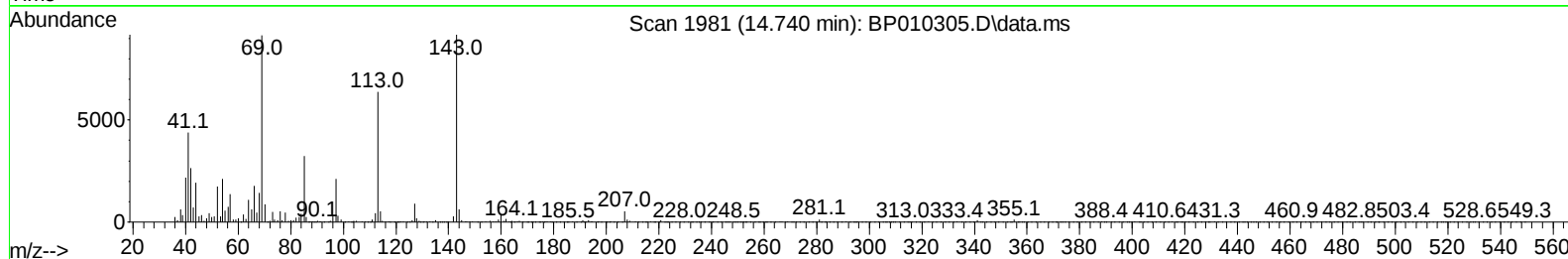
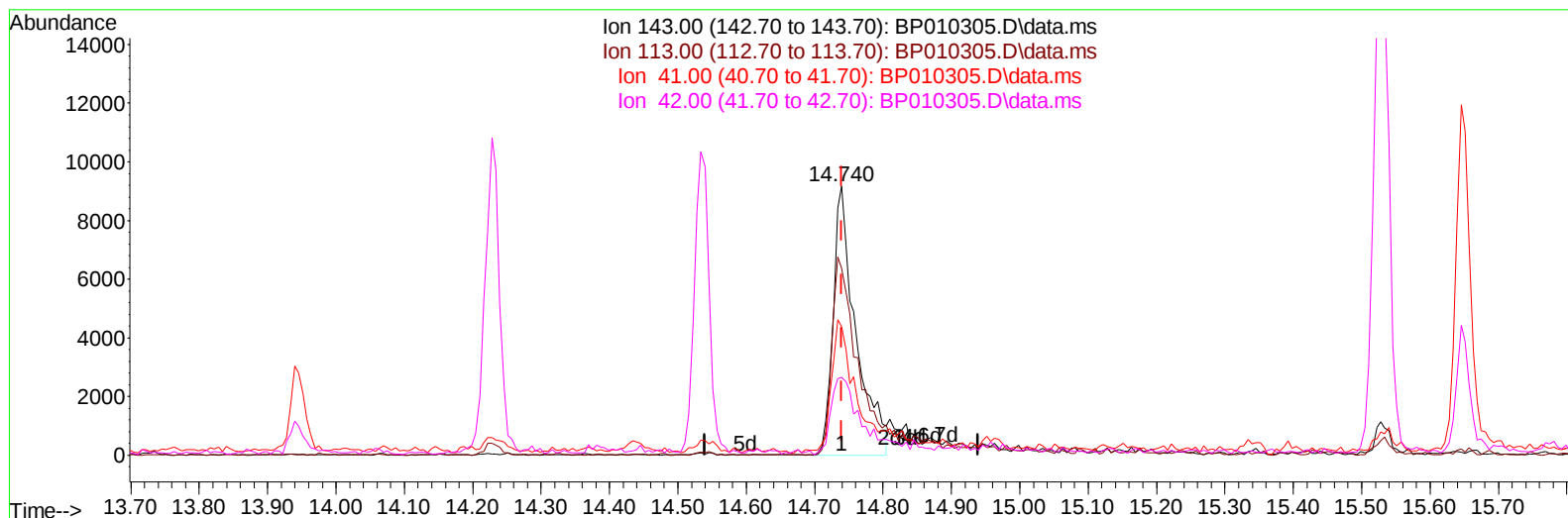
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TIC: BP010305.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.000) 3.36 ng/u1

response 21170

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	69.36#
41.00	23.20	47.66#
42.00	18.00	28.98#

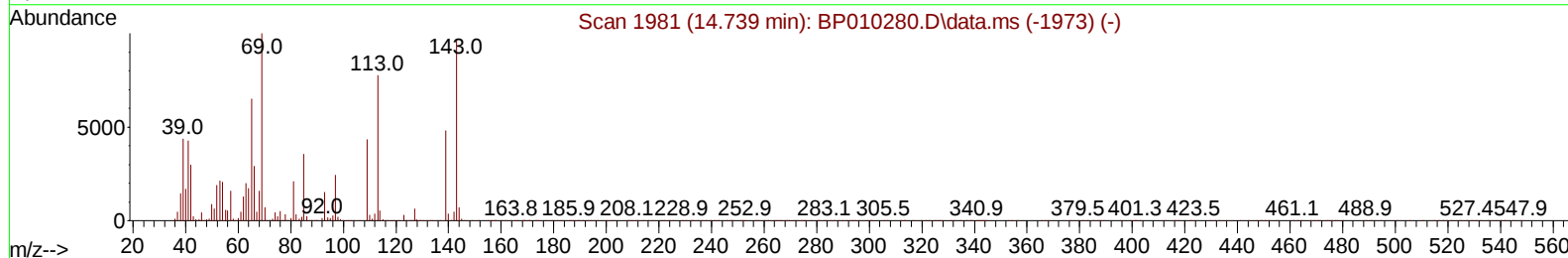
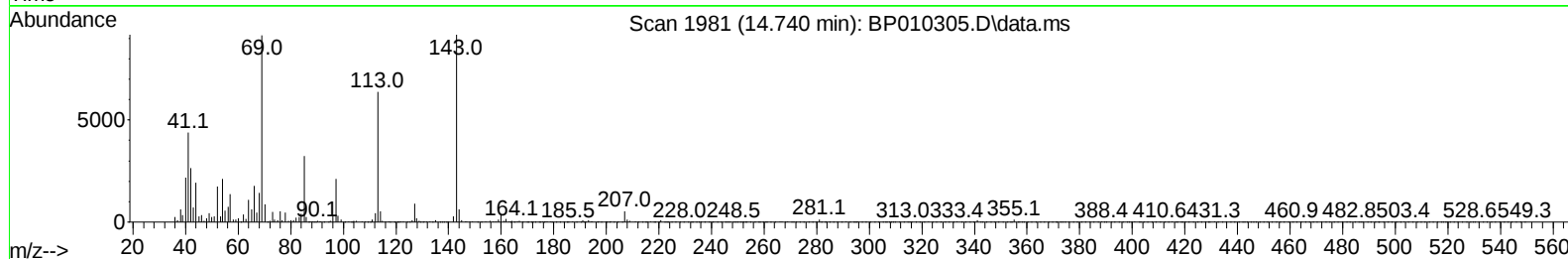
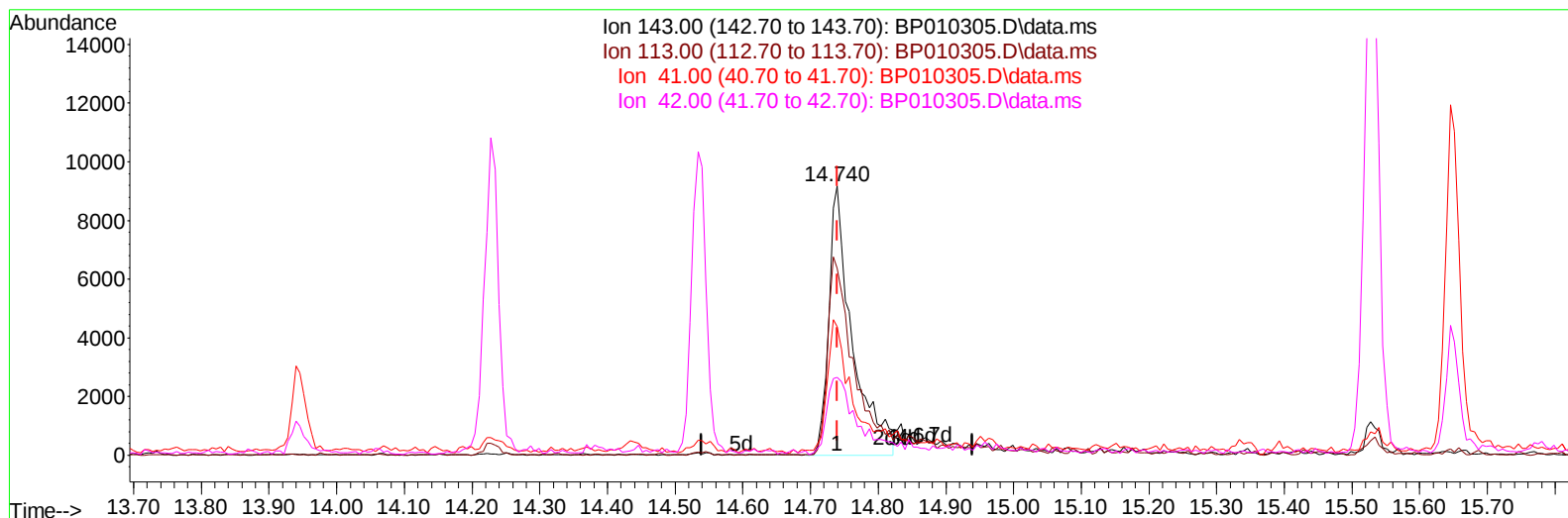
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Instrument :
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TIC: BP010305.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.000) 3.52 ng/ul m

response 22223

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	69.36#
41.00	23.20	47.66#
42.00	18.00	28.98#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	137681	20.000	ng/ul	0.00
20) Naphthalene-d8	10.705	136	579508	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164	412710	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.287	188	932771	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.369	240	936821	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	906061	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	16235m	5.009	ng/uL	0.00
4) Pyridine-d5	3.775	84	45975	5.098	ng/ul	0.00
7) Phenol-d5	7.069	99	64236	5.023	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	231871	29.311	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	206993	21.979	ng/ul	0.00
15) 4-Methylphenol-d8	8.616	113	131042	12.465	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	141800	30.517	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	145061	27.381	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	247038	25.619	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131	210287	14.772	ng/ul	0.00
46) Dimethylphthalate-d6	13.946	166	1021289	31.062	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	1118716	28.538	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	22223m	3.524	ng/ul	0.00
60) Fluorene-d10	15.528	176	890793	32.096	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	171548	26.857	ng/ul	0.00
73) Anthracene-d10	17.387	188	1595264	34.924	ng/ul	0.00
81) Pyrene-d10	19.616	212	1920961	37.106	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.645	264	1696796	35.116	ng/ul	-0.01

Target Compounds Qvalue

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

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