

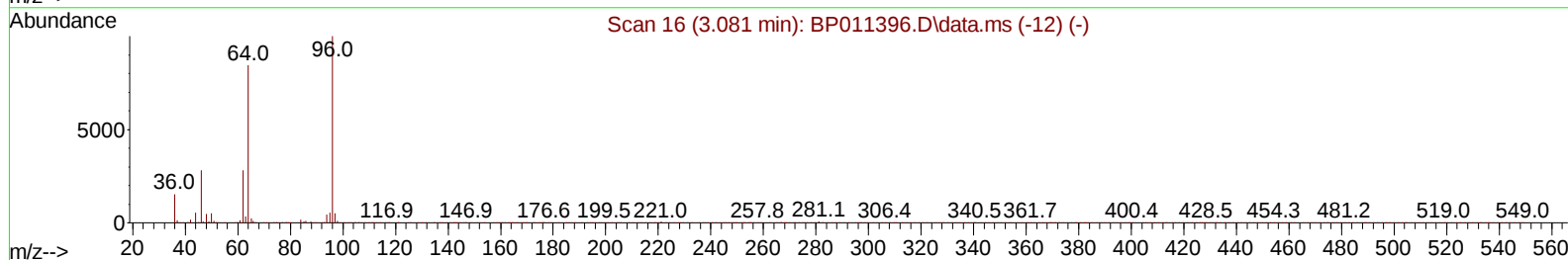
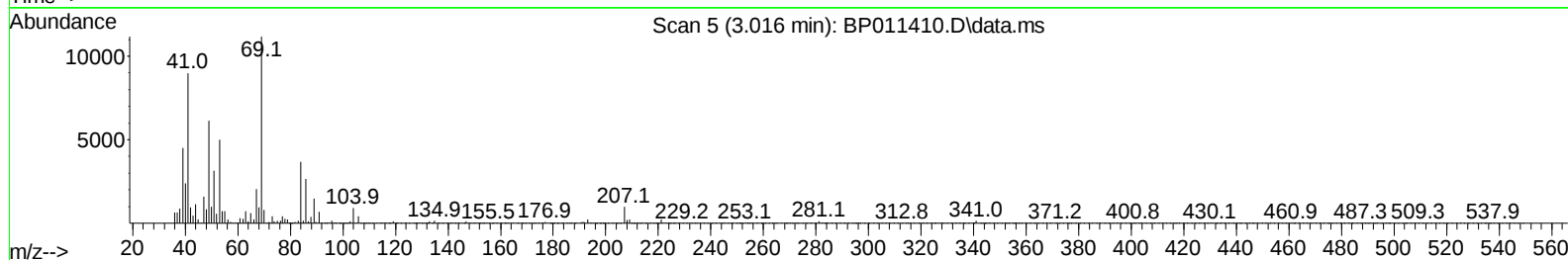
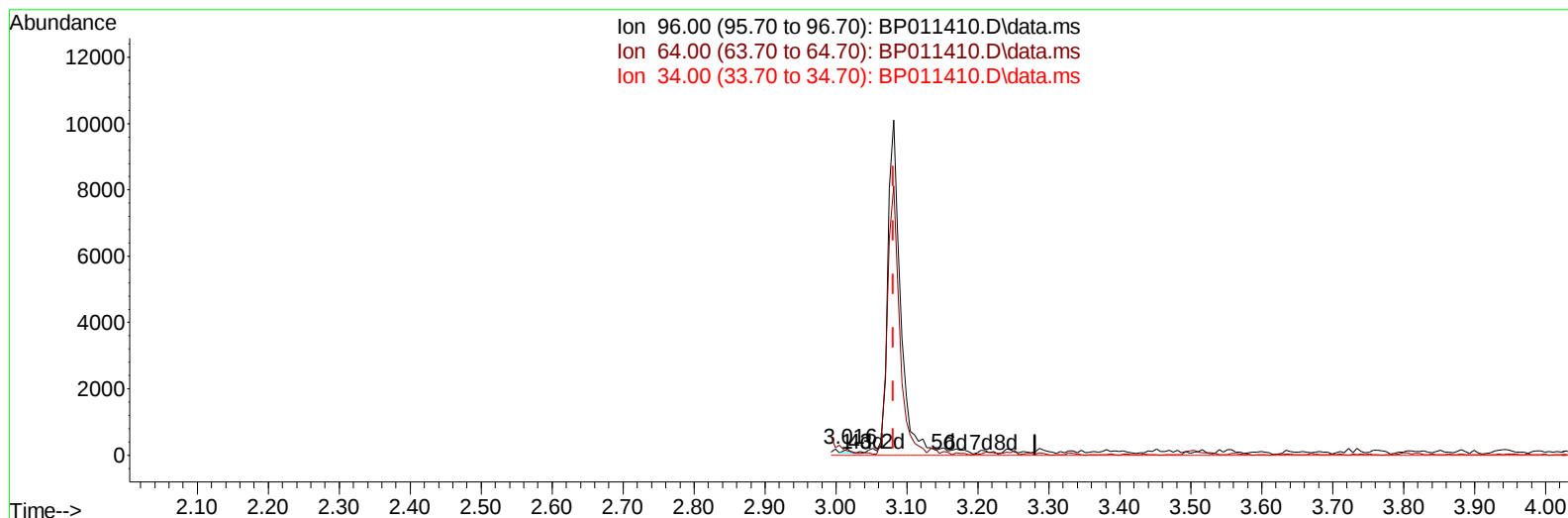
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
 Data File : BP011410.D
 Acq On : 12 Aug 2022 04:51
 Operator : CG/JU
 Sample : N4129-02
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGBW2

Manual IntegrationsAPPROVED

Quant Time: Aug 12 05:35:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 11 23:05:53 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/12/2022
 Supervised By :Sohil Jodhani 08/13/2022



TIC: BP011410.D\data.ms

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

3.016min (-0.065) 0.03 ng/uL

response 68

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	64.74
34.00	0.00	0.00
0.00	0.00	0.00

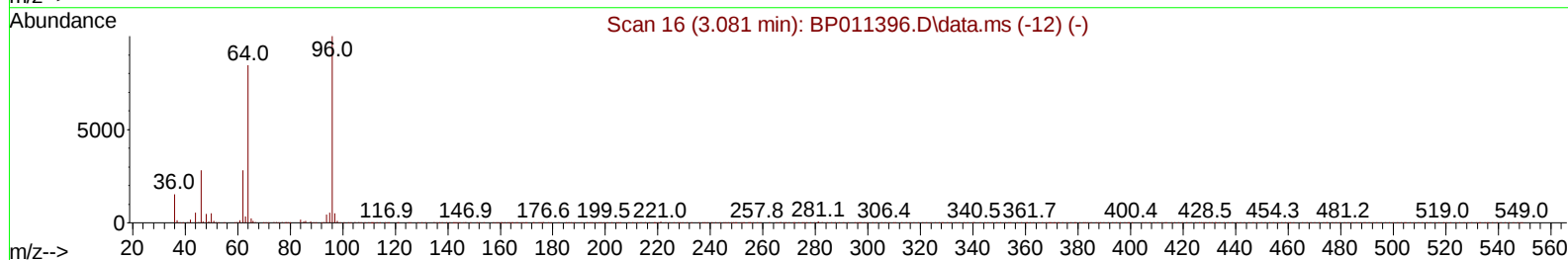
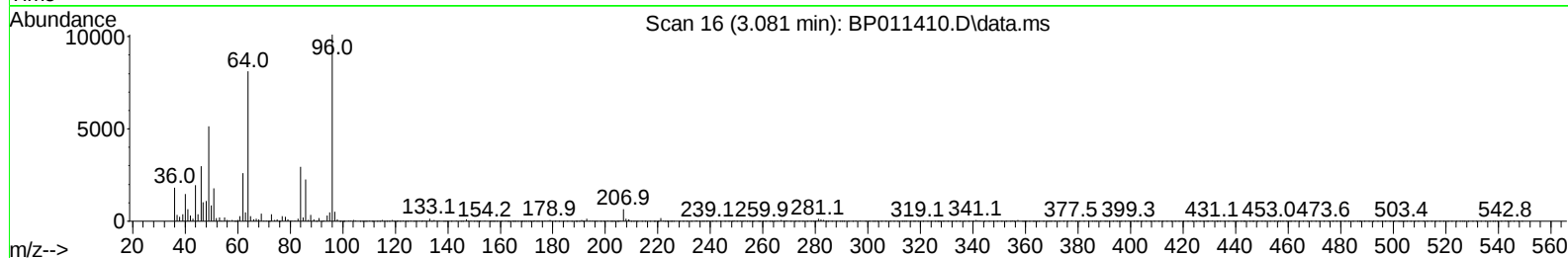
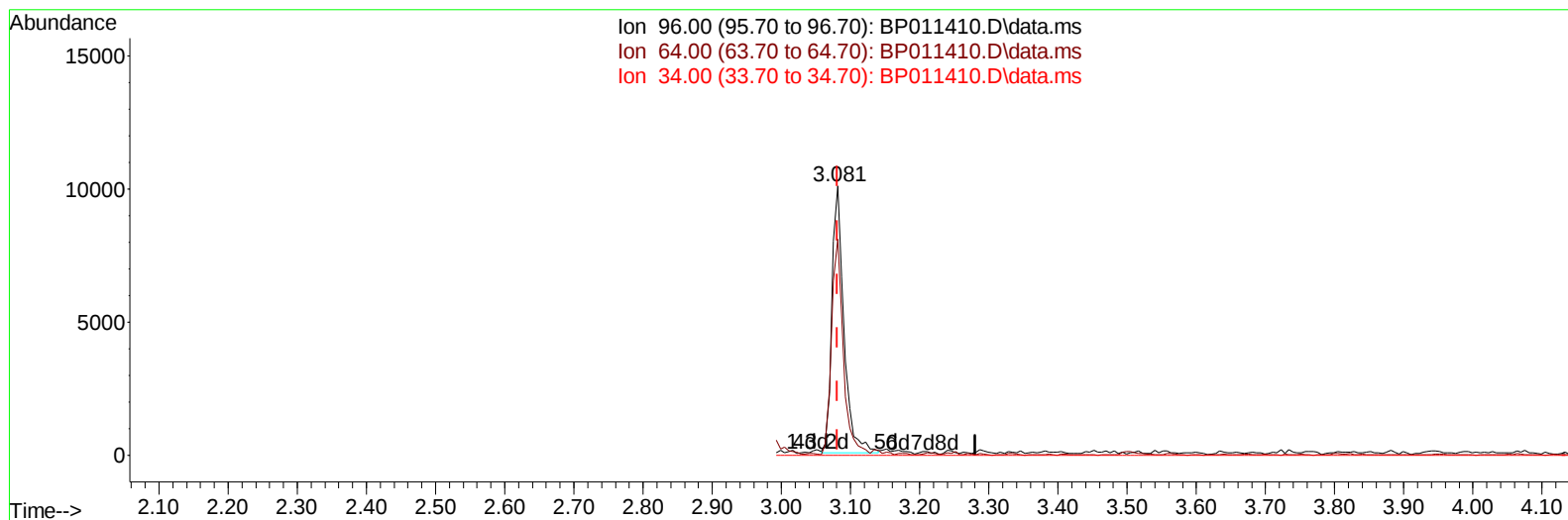
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(3) 1,4-Dioxane-d8 (S)

3.081min (-0.000) 4.49 ng/uL m

response 12106

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	80.28#
34.00	0.00	0.00
0.00	0.00	0.00

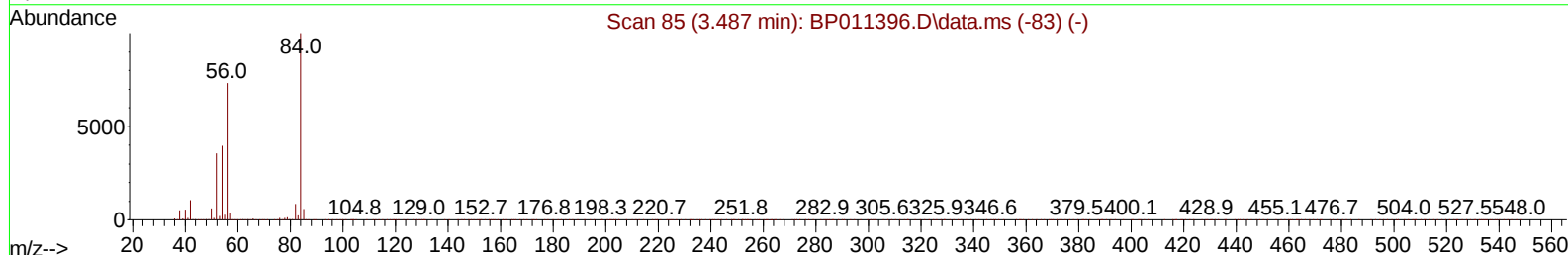
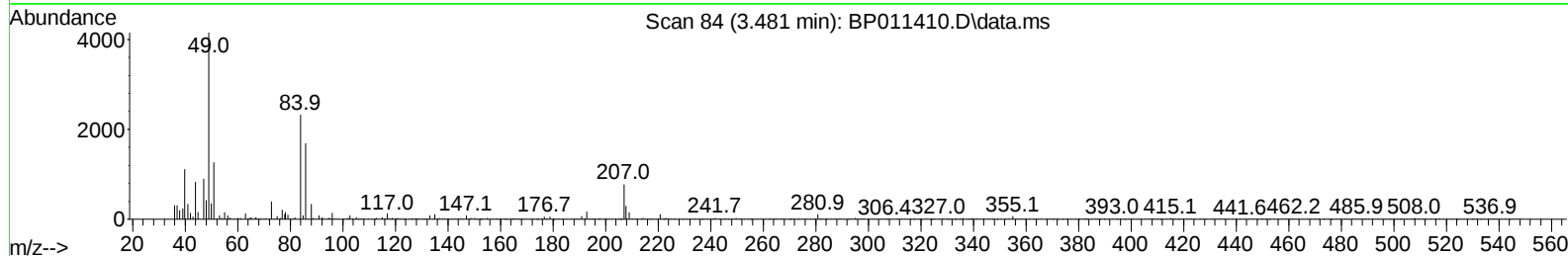
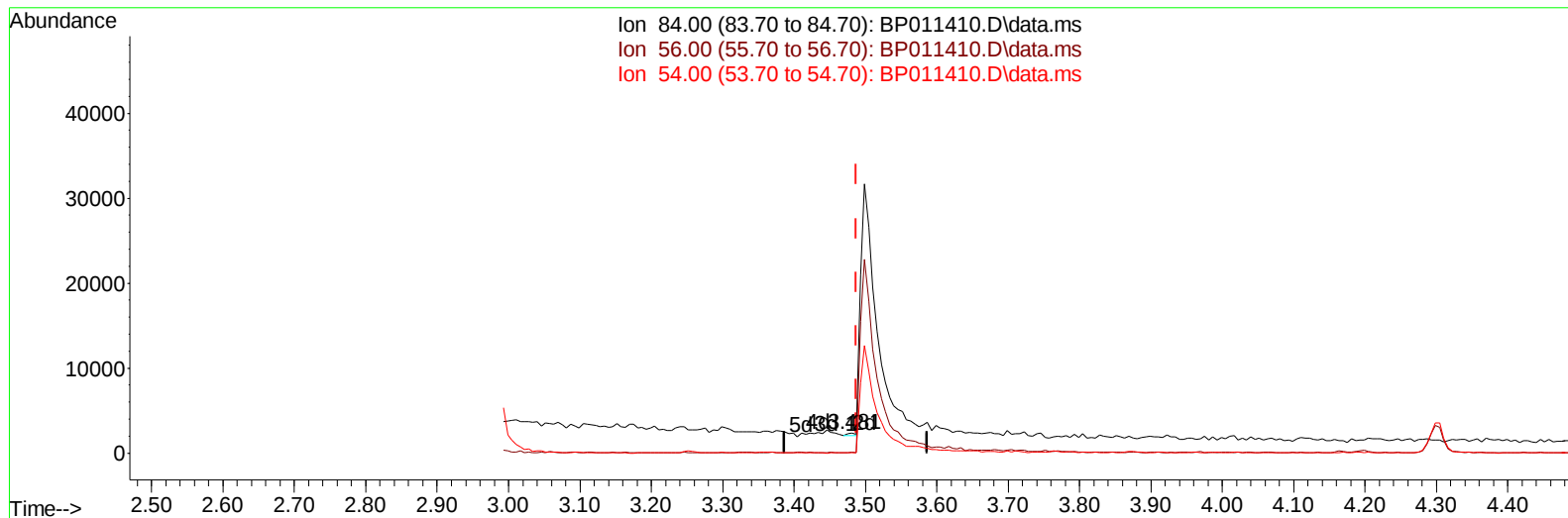
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TIC: BP011410.D\data.ms

(4) Pyridine-d5 (S)

3.481min (-0.006) 0.03 ng/u1

response 223

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.60	3.78#
54.00	33.20	1.16#
0.00	0.00	0.00

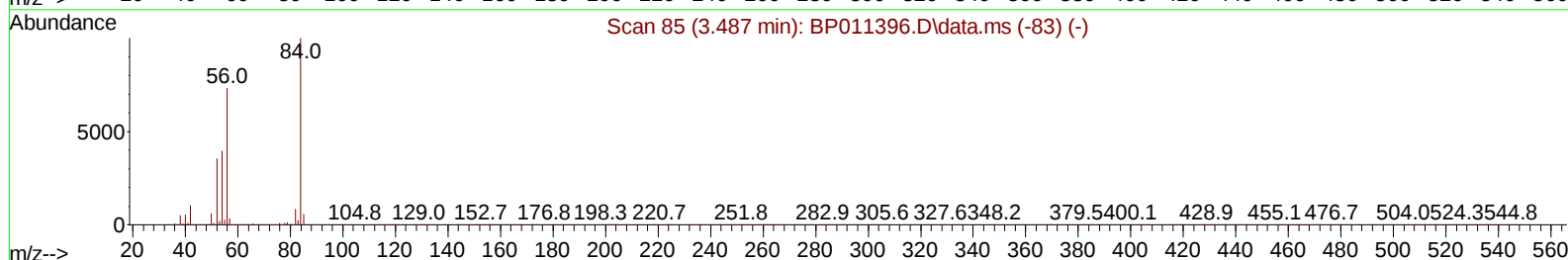
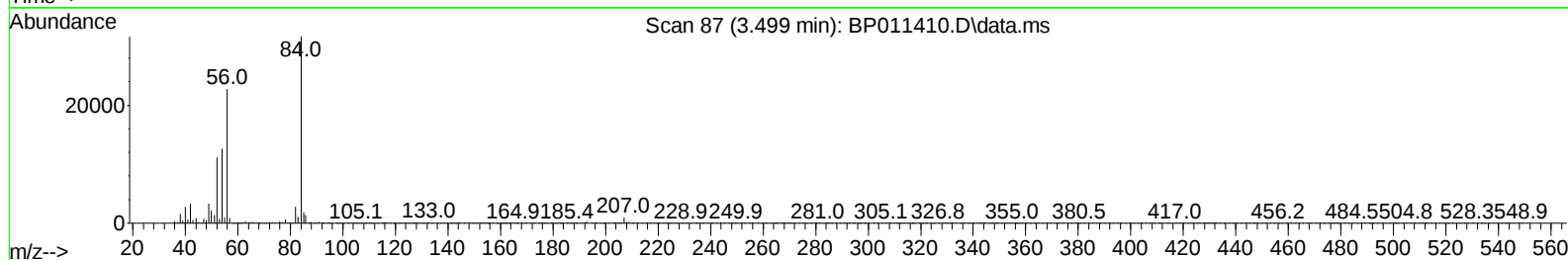
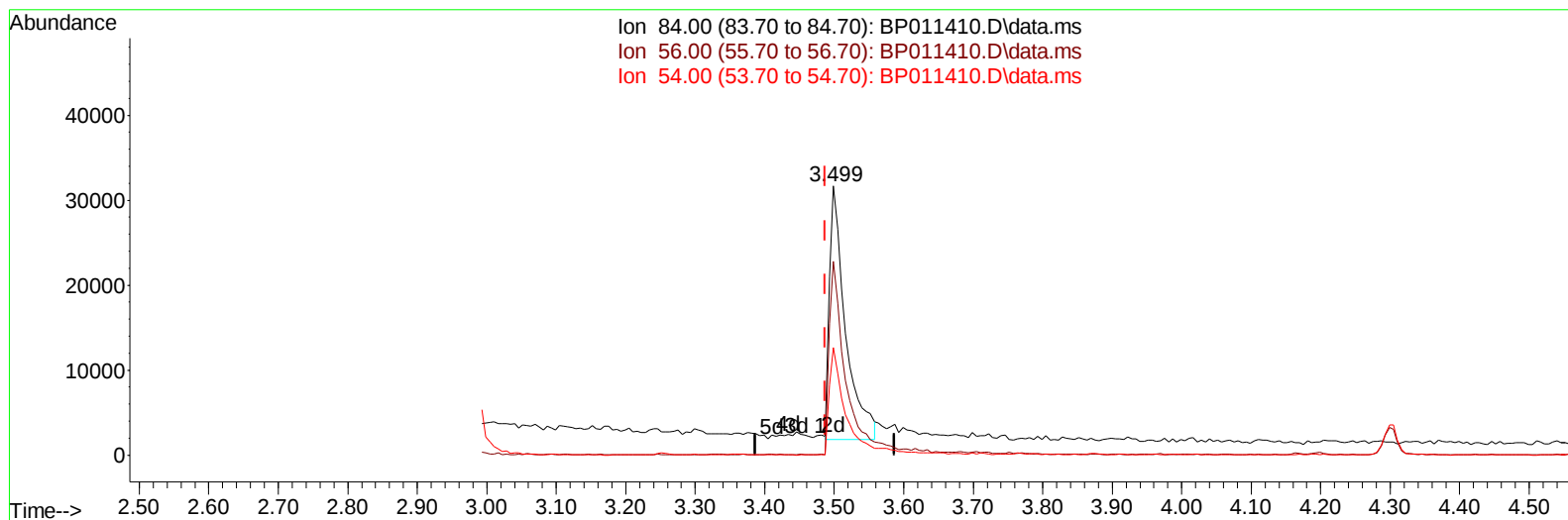
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 ALS Vial : 33 Sample Multiplier: 1

Instrument :
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TIC: BP011410.D\data.ms

(4) Pyridine-d5 (S)

3.499min (+ 0.012) 7.03 ng/u1 m

response 47871

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.60	71.97
54.00	33.20	40.00#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	82744	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	349651	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.251	164	197842	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.022	188	417276	20.000	ng/ul	0.00
79) Chrysene-d12	21.151	240	404412	20.000	ng/ul	0.00
88) Perylene-d12	23.457	264	424242	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	12106m	4.485	ng/uL	0.00
4) Pyridine-d5	3.499	84	47871m	7.031	ng/ul	0.01
7) Phenol-d5	6.763	99	5671	0.774	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	159588	30.737	ng/ul	0.00
11) 2-Chlorophenol-d4	7.110	132	76193	13.261	ng/ul	0.00
15) 4-Methylphenol-d8	8.299	113	19008	3.306	ng/ul	0.00
21) Nitrobenzene-d5	8.740	128	90287	37.942	ng/ul	0.00
24) 2-Nitrophenol-d4	9.457	143	72700	33.369	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	114880	26.111	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	1109	0.144	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	559295	40.743	ng/ul	0.00
49) Acenaphthylene-d8	13.939	160	291884	17.664	ng/ul	0.00
54) 4-Nitrophenol-d4	14.492	143	7942	3.168	ng/ul	0.01
60) Fluorene-d10	15.251	176	389090	33.144	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.386	200	53981	27.851	ng/ul	0.00
73) Anthracene-d10	17.122	188	496505	26.472	ng/ul	0.00
81) Pyrene-d10	19.386	212	676897	30.381	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.309	264	579129	27.069	ng/ul	0.00

Target Compounds Qvalue

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

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