

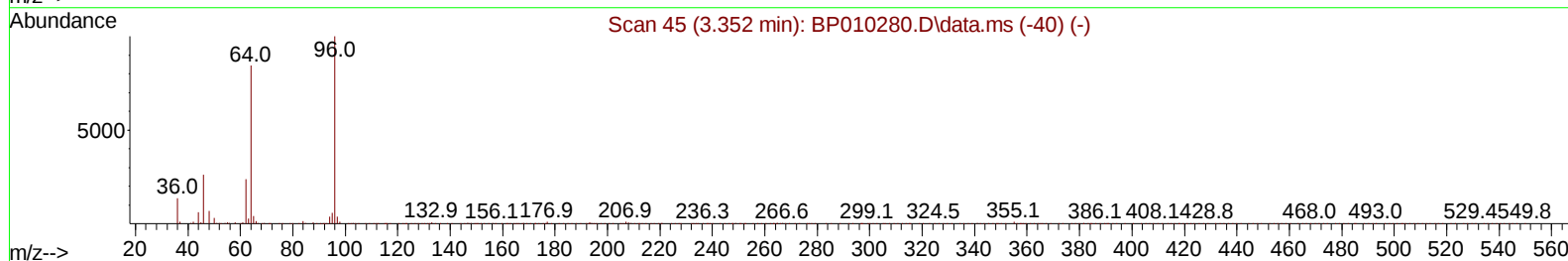
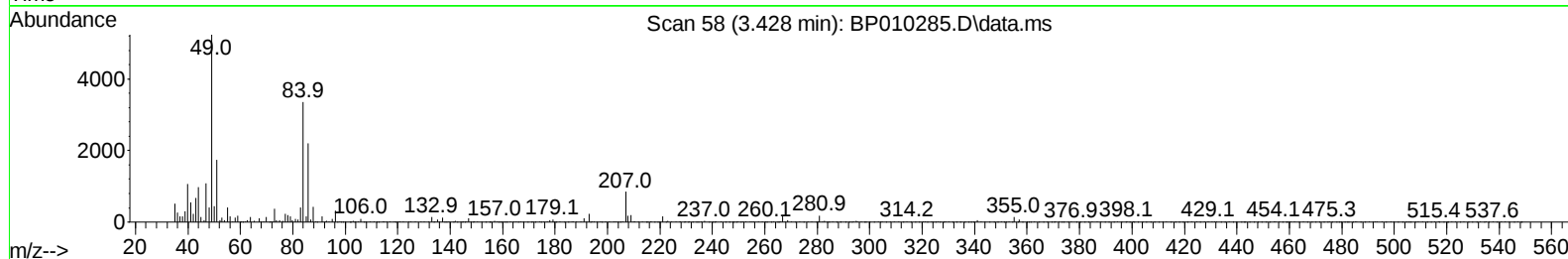
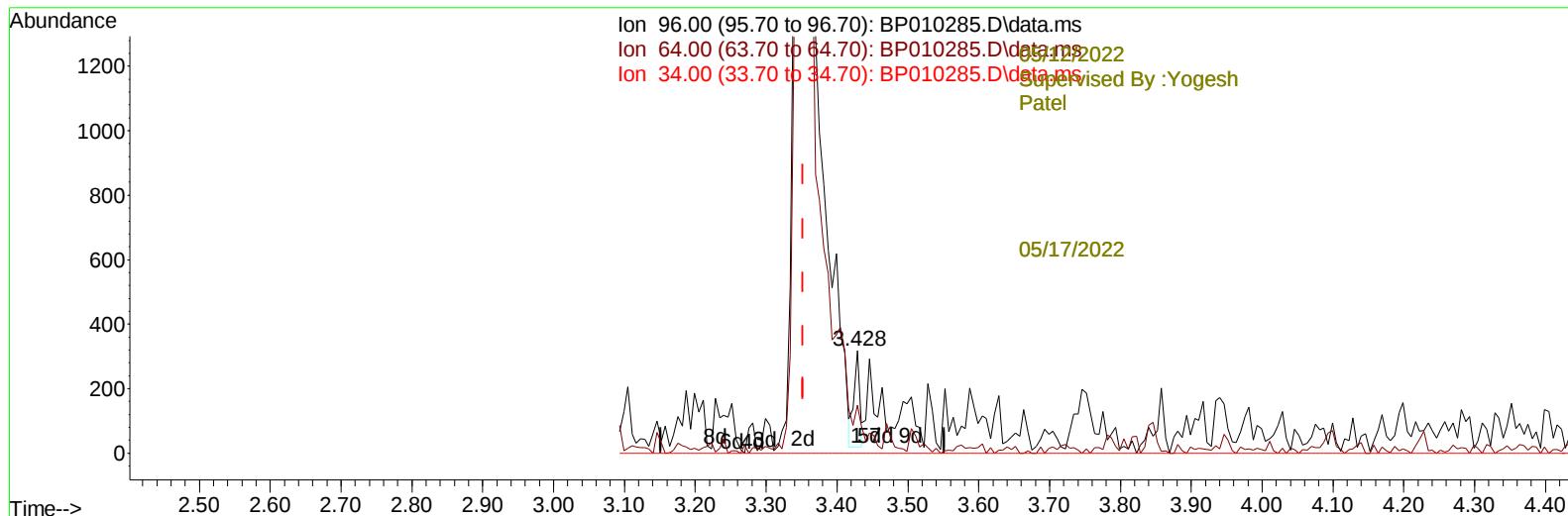
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010285.D
 Acq On : 11 May 2022 12:35
 Operator : CG/JU
 Sample : N2632-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW4A4

Manual IntegrationsAPPROVED

Quant Time: May 12 01:12:28 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

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



TIC: BP010285.D\data.ms

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

3.428min (+ 0.077) 0.09 ng/uL

response 174

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	46.71
34.00	0.00	0.00
0.00	0.00	0.00

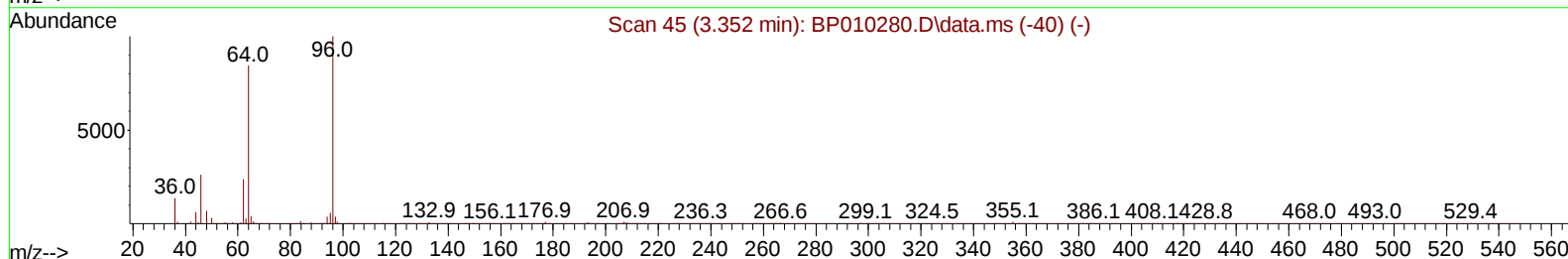
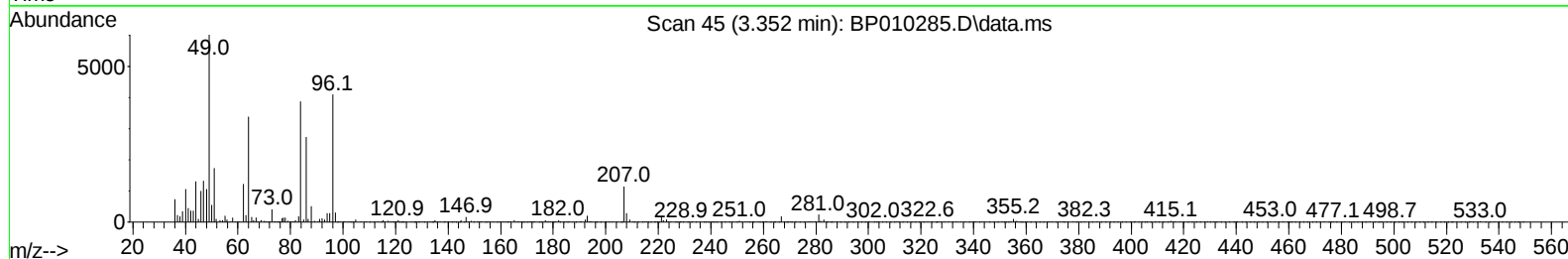
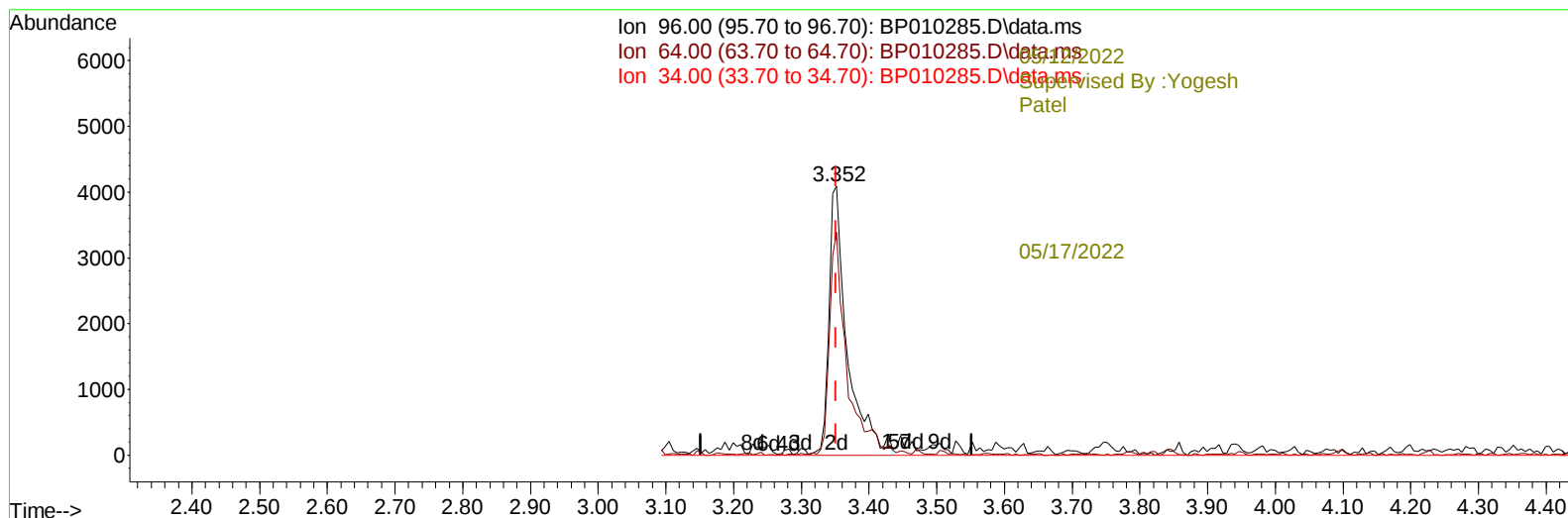
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TIC: BP010285.D\data.ms

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

3.352min (+ 0.000) 4.17 ng/uL m

response 7659

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	82.91#
34.00	0.00	0.00
0.00	0.00	0.00

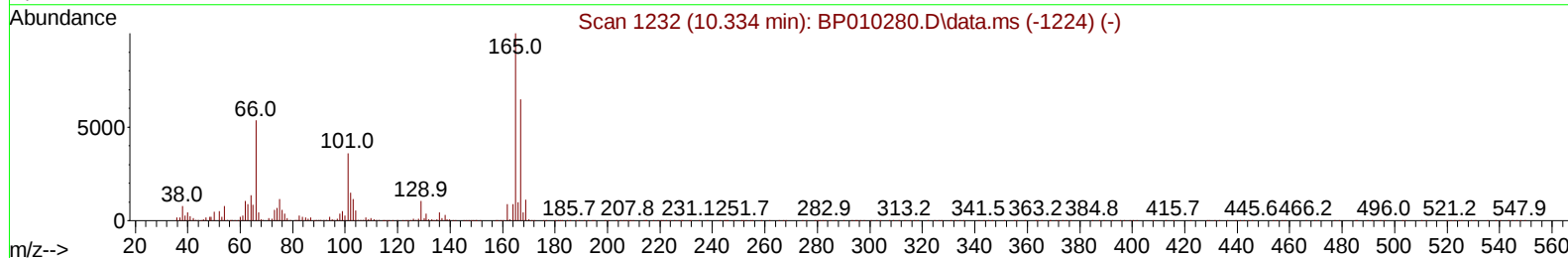
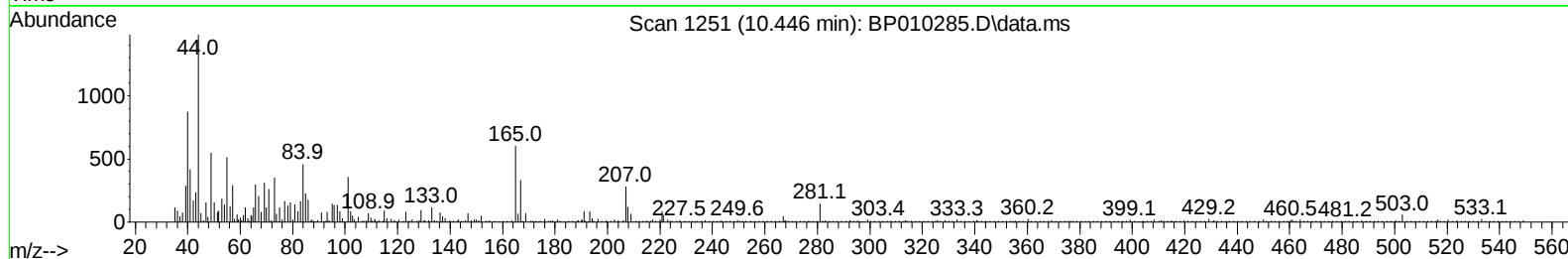
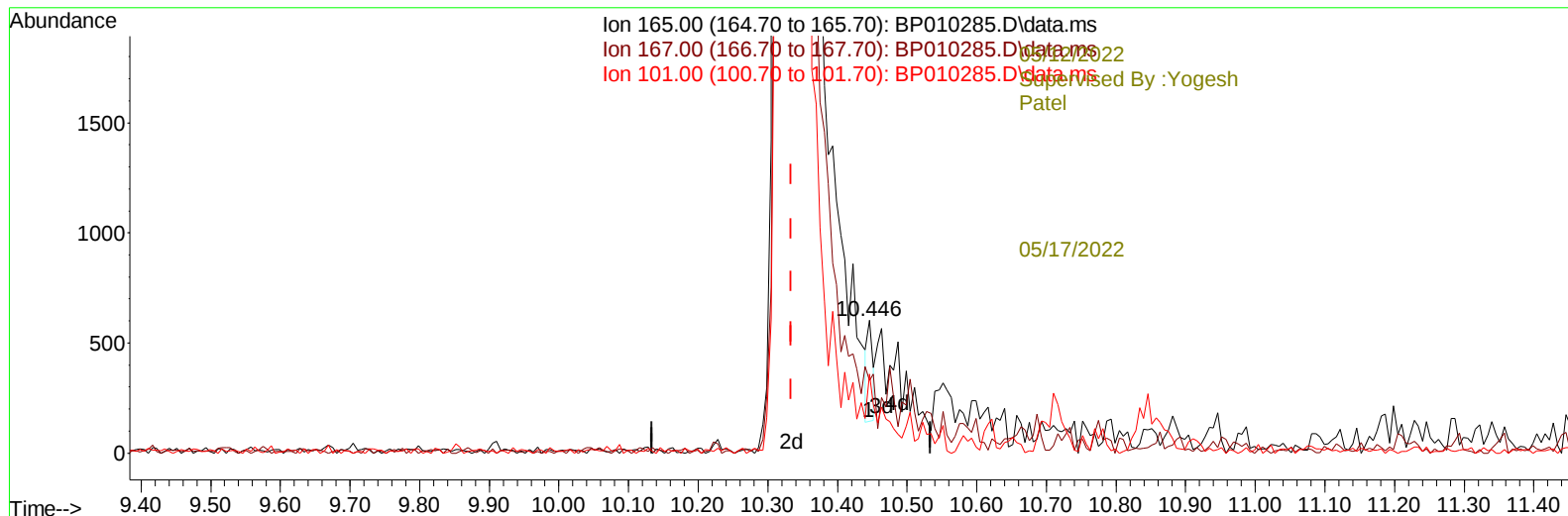
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TIC: BP010285.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.446min (+ 0.112) 0.04 ng/ul

response 247

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	54.89
101.00	49.80	59.54
0.00	0.00	0.00

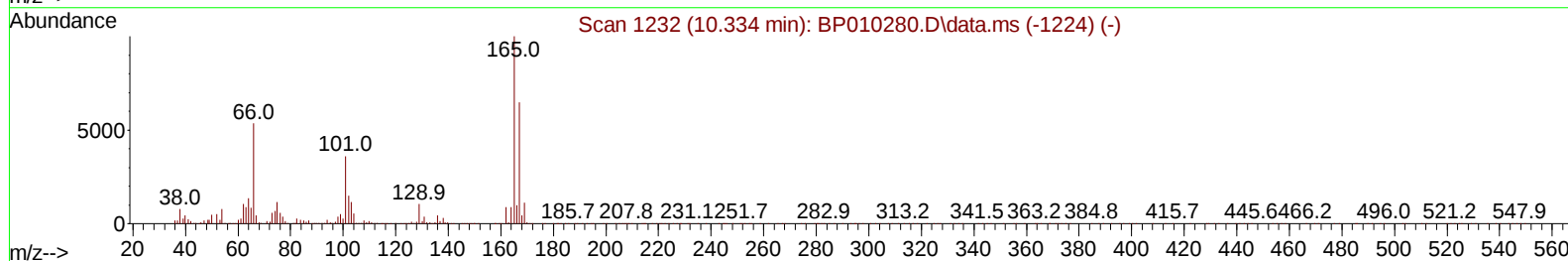
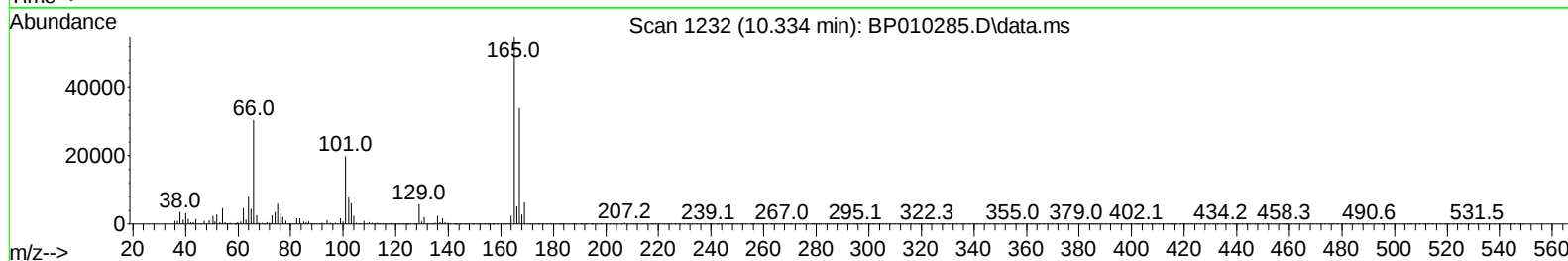
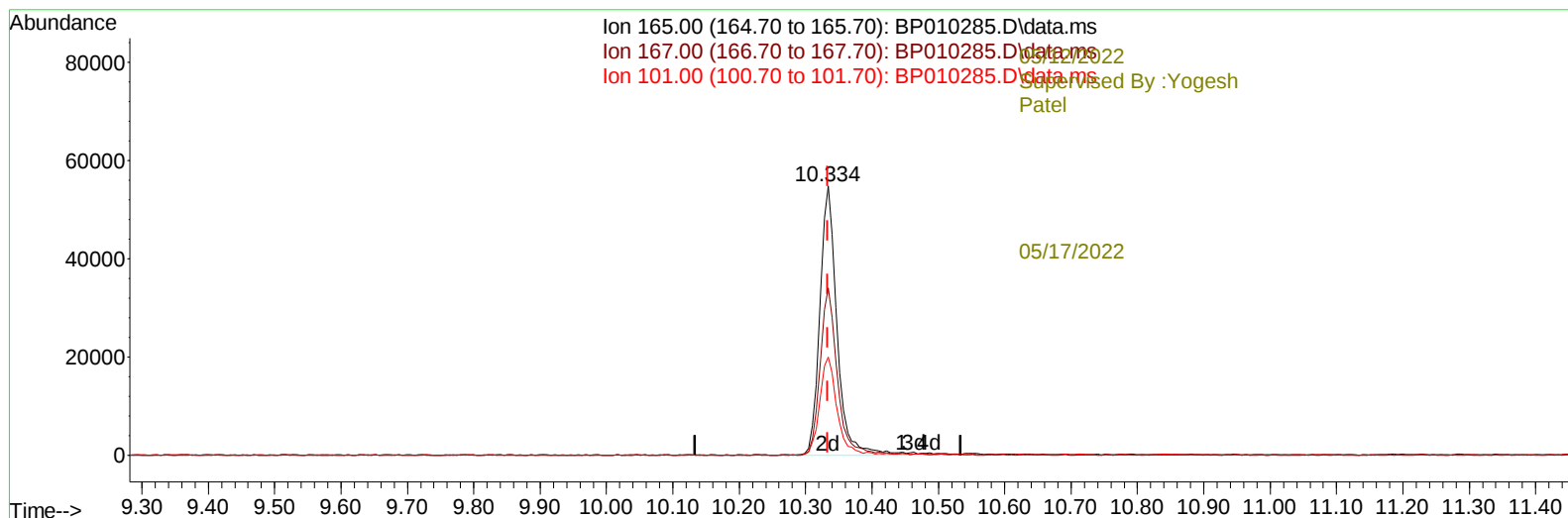
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TIC: BP010285.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.334min (+ 0.000) 17.39 ng/ul m

response 99115

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	62.18
101.00	49.80	36.35#
0.00	0.00	0.00

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
-----							05/12/2022
Internal Standards							Supervised By :Yogesh Patel
1) 1,4-Dichlorobenzene-d4	7.911	152		78087	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136		342538	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.540	164		244525	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.287	188		547234	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.375	240		572339	20.000	ng/ul	0.00
88) Perylene-d12	23.804	264		615772	20.000	ng/ul	-0.01
-----							05/17/2022
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.352	96		7659m	4.167	ng/uL	0.00
4) Pyridine-d5	3.787	84		23221	4.540	ng/ul	0.02
7) Phenol-d5	7.069	99		127589	17.592	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67		94145	20.983	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132		104872	19.634	ng/ul	0.00
15) 4-Methylphenol-d8	8.616	113		107817	18.082	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128		54740	19.931	ng/ul	0.00
24) 2-Nitrophenol-d4	9.793	143		55821	17.826	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.334	165		99115m	17.390	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131		137129	16.297	ng/ul	0.00
46) Dimethylphthalate-d6	13.946	166		385616	19.795	ng/ul	0.00
49) Acenaphthylene-d8	14.234	160		427288	18.397	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143		33693	9.017	ng/ul	0.00
60) Fluorene-d10	15.534	176		340623	20.714	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200		34123	9.106	ng/ul	0.00
73) Anthracene-d10	17.387	188		545051	20.339	ng/ul	0.00
81) Pyrene-d10	19.622	212		701414	22.177	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.645	264		705604	21.487	ng/ul	-0.01

Target Compounds						Qvalue	
16) Acetophenone	8.734	105		9612	1.031	ng/ul#	66

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

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