

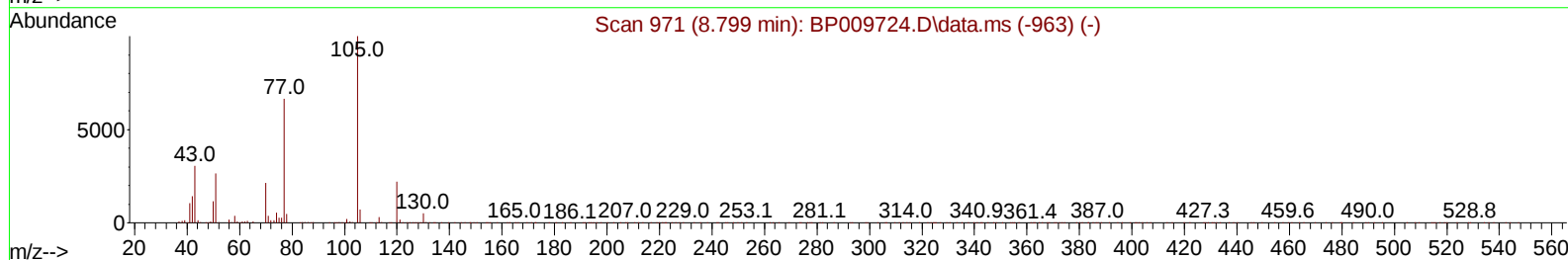
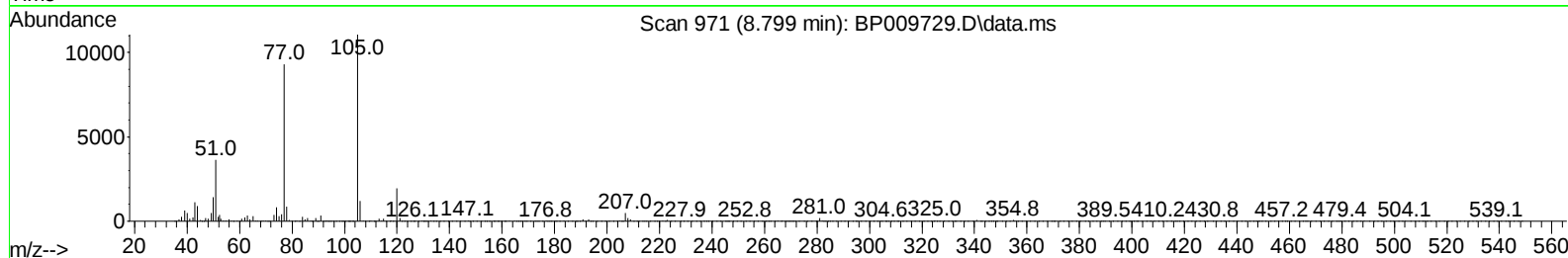
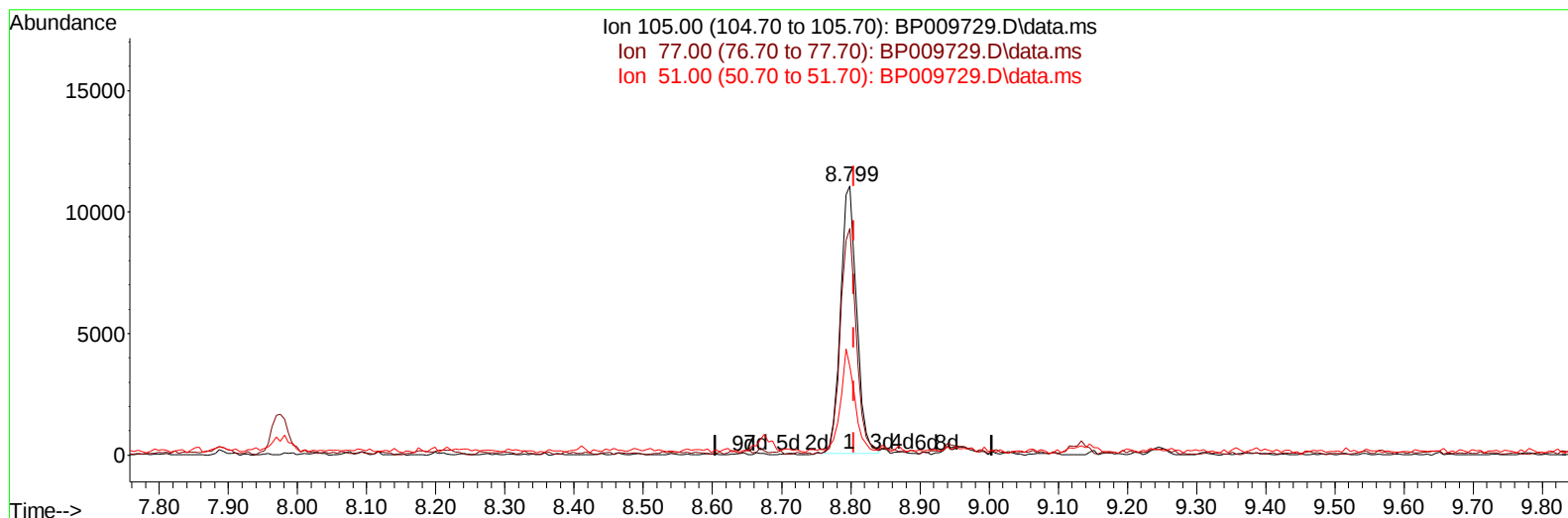
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
 Data File : BP009729.D
 Acq On : 08 Apr 2022 22:09
 Operator : CG/JU
 Sample : N2182-04
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0R13

Manual IntegrationsAPPROVED

Quant Time: Apr 09 01:46:43 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 07 19:16:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/11/2022
 Supervised By :Yogesh Patel 04/12/2022



TIC: BP009729.D\data.ms

(16) Acetophenone

8.799min (-0.006) 2.20 ng/u1

response 17899

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	66.80	84.24#
51.00	17.30	32.86#
0.00	0.00	0.00

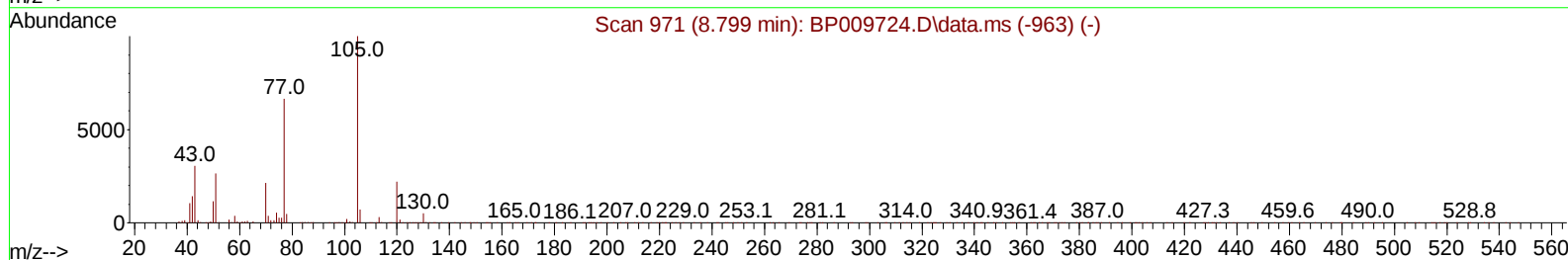
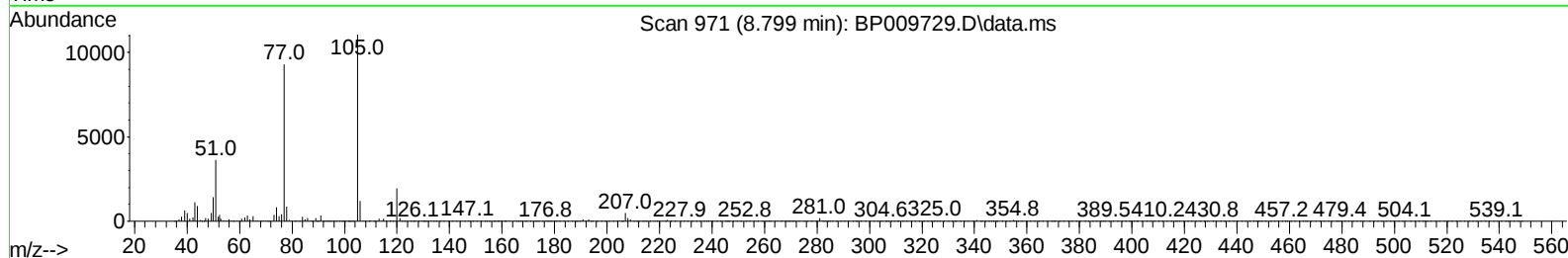
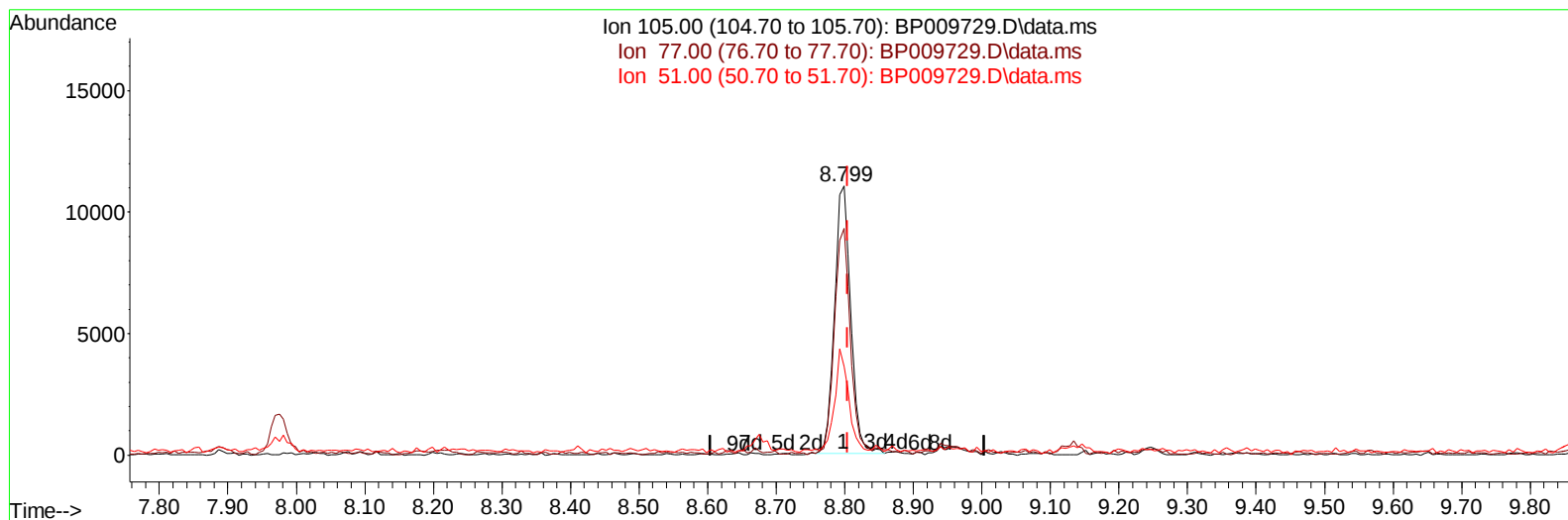
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TIC: BP009729.D\data.ms

(16) Acetophenone

8.799min (-0.006) 2.22 ng/ul m

response 18064

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	66.80	84.24#
51.00	17.30	32.86#
0.00	0.00	0.00

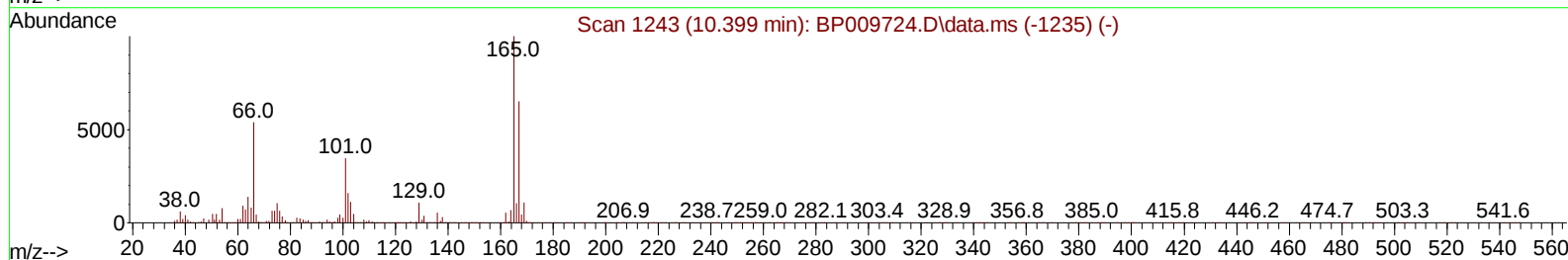
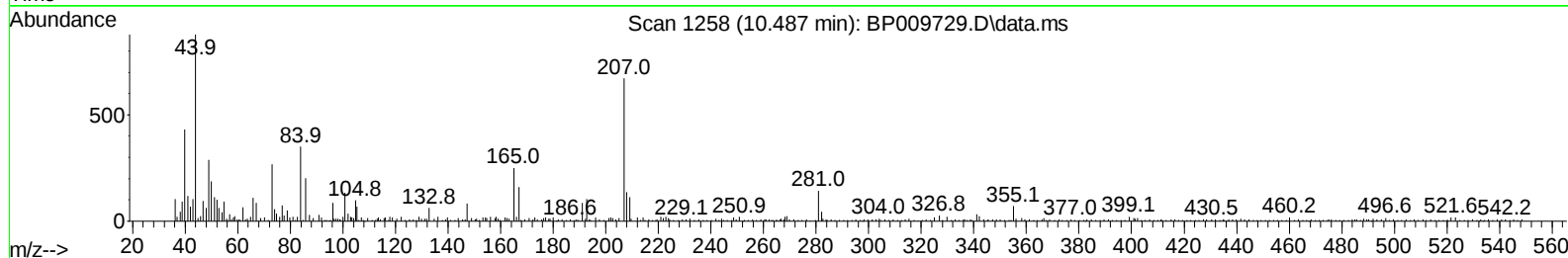
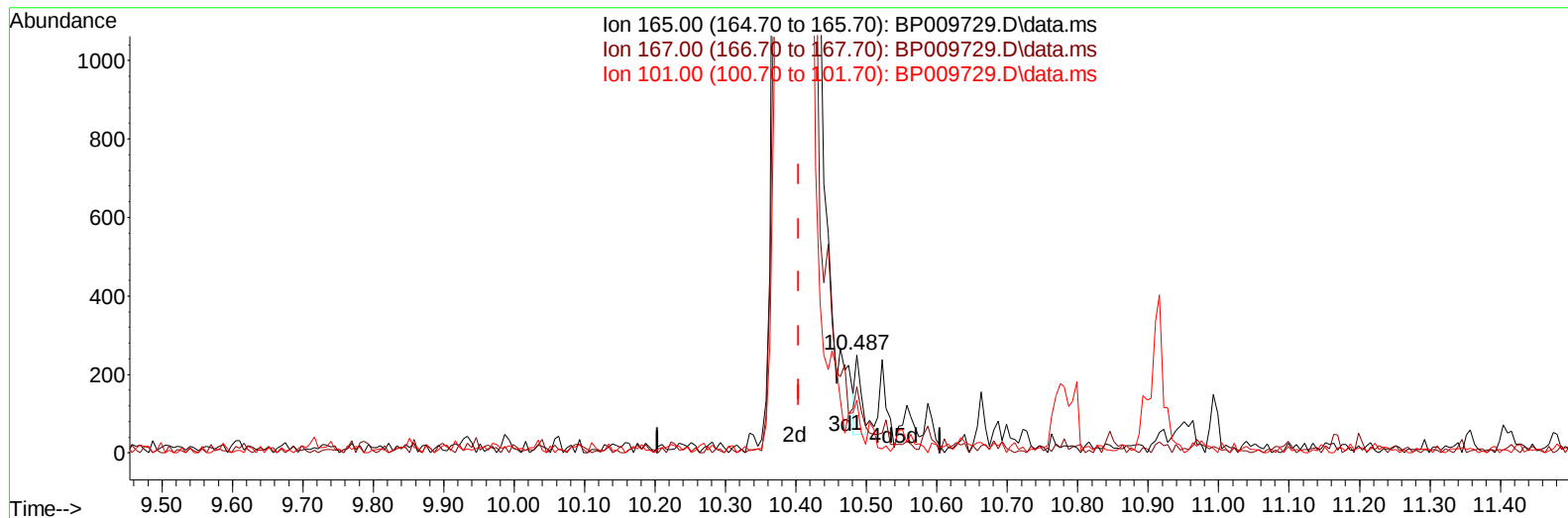
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TIC: BP009729.D\data.ms

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

10.487min (+ 0.082) 0.02 ng/u1

response 115

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	67.60
101.00	49.80	54.00
0.00	0.00	0.00

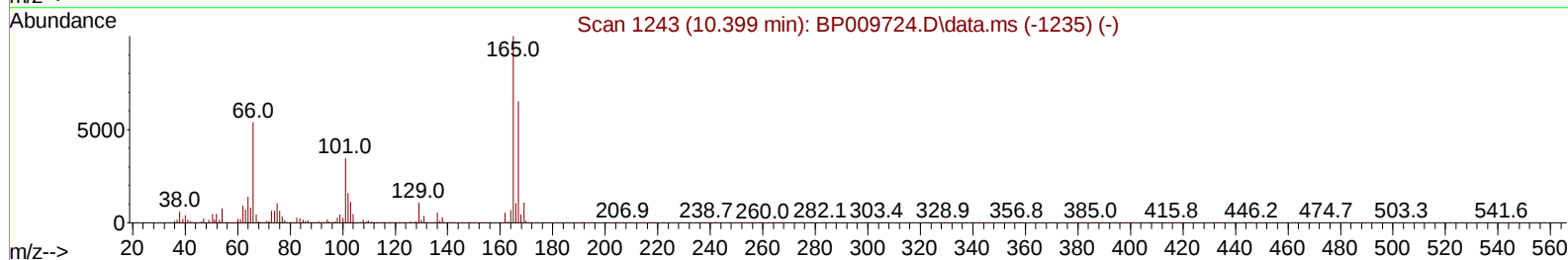
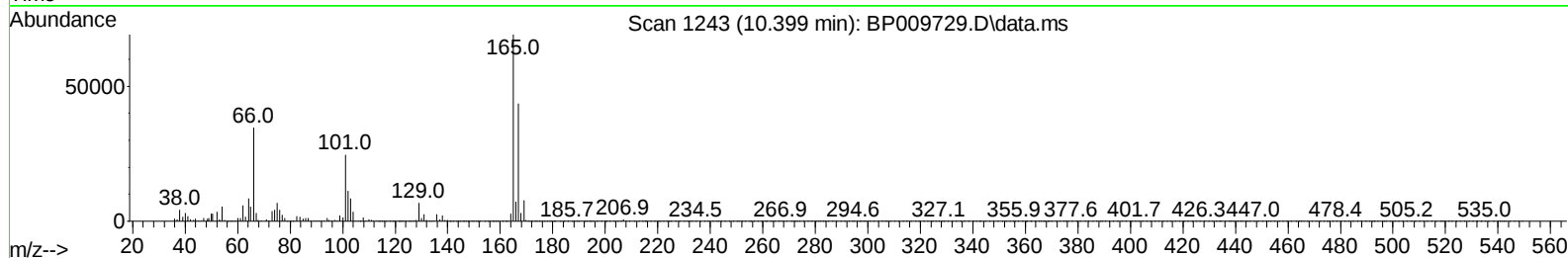
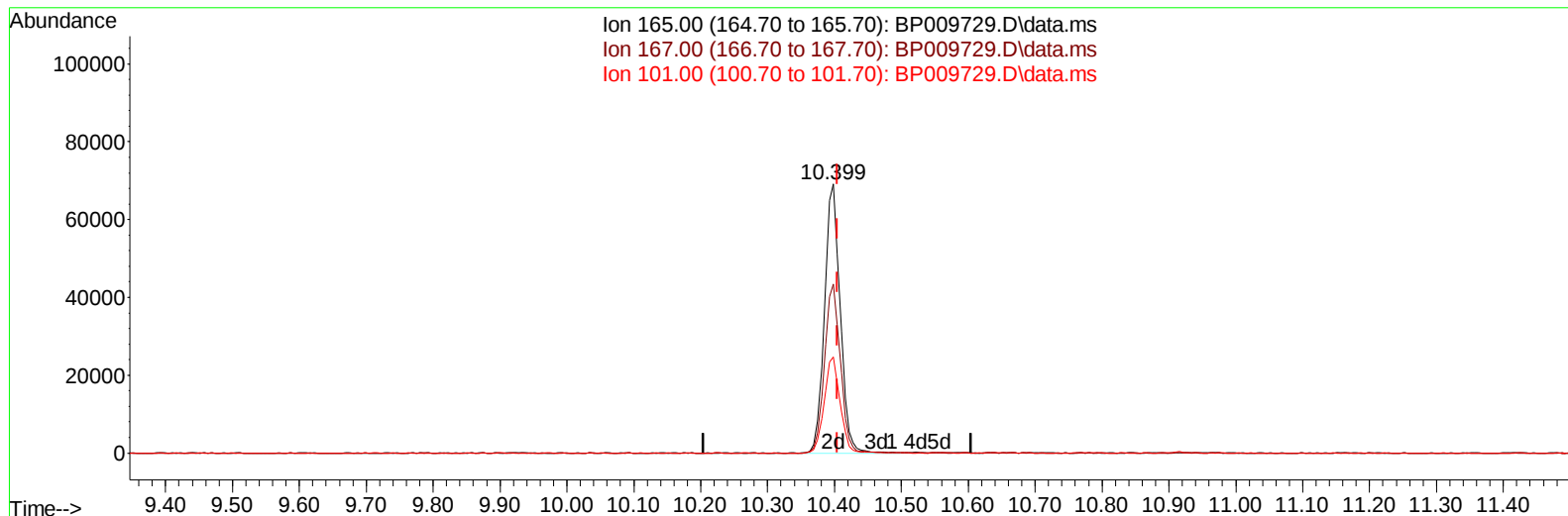
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TIC: BP009729.D\data.ms

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

10.399min (-0.006) 21.10 ng/ul m

response 113072

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	62.93
101.00	49.80	35.66#
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	7.975	152	70607	20.000	ng/ul	0.00
20) Naphthalene-d8	10.781	136	323940	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.604	164	240061	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.357	188	541904	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.433	240	593449	20.000	ng/ul	-0.01
88) Perylene-d12	23.910	264	625402	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	6273	3.965	ng/uL	0.00
4) Pyridine-d5	3.811	84	37578	8.615	ng/ul	0.00
7) Phenol-d5	7.122	99	129432	20.880	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	87503	23.711	ng/ul	0.00
11) 2-Chlorophenol-d4	7.499	132	102099	21.835	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	111523	21.551	ng/ul	0.00
21) Nitrobenzene-d5	9.128	128	55568	23.787	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.857	143	60133	23.479	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	113072m	21.098	ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131	155407	20.360	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	497205	26.367	ng/ul	-0.01
49) Acenaphthylene-d8	14.298	160	506269	22.662	ng/ul	0.00
54) 4-Nitrophenol-d4	14.781	143	93685	27.395	ng/ul	0.00
60) Fluorene-d10	15.598	176	420208	26.489	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	73628	22.935	ng/ul	0.00
73) Anthracene-d10	17.451	188	762965	29.406	ng/ul	-0.01
81) Pyrene-d10	19.680	212	1038214	32.032	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.751	264	1079904	33.158	ng/ul	-0.01
Target Compounds						
						Qvalue
16) Acetophenone	8.799	105	18064m	2.216	ng/ul	

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

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