

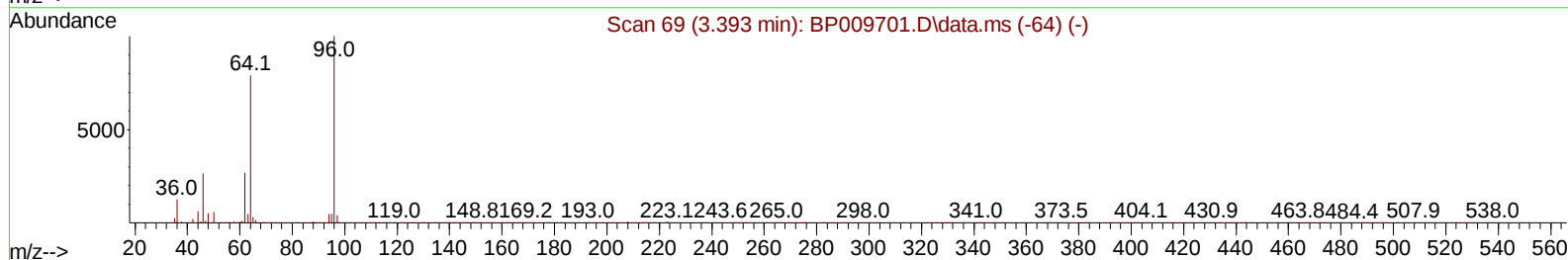
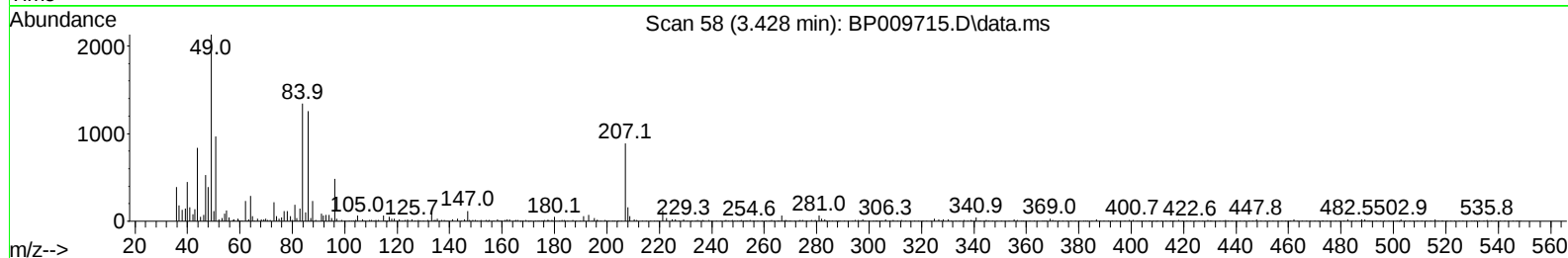
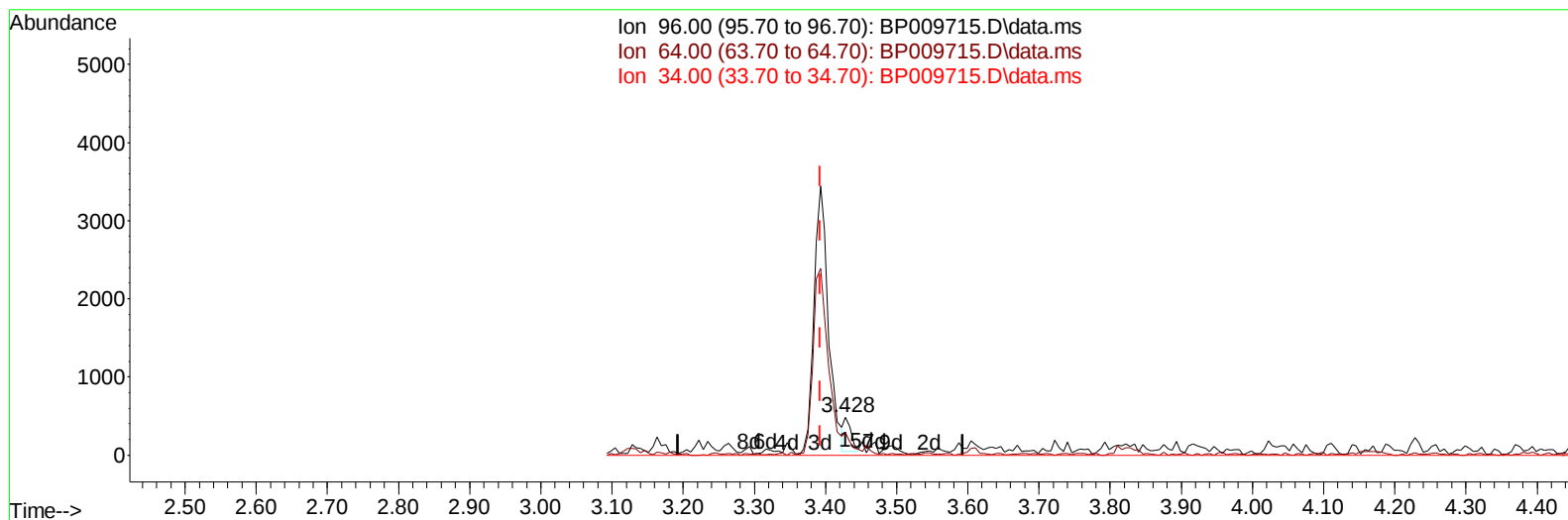
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
 Data File : BP009715.D
 Acq On : 08 Apr 2022 13:41
 Operator : CG/JU
 Sample : N2183-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0R30

Manual IntegrationsAPPROVED

Quant Time: Apr 08 22:27:03 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: BP009715.D\data.ms

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

3.428min (+ 0.035) 0.15 ng/uL

response 311

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	59.59
34.00	0.00	0.00
0.00	0.00	0.00

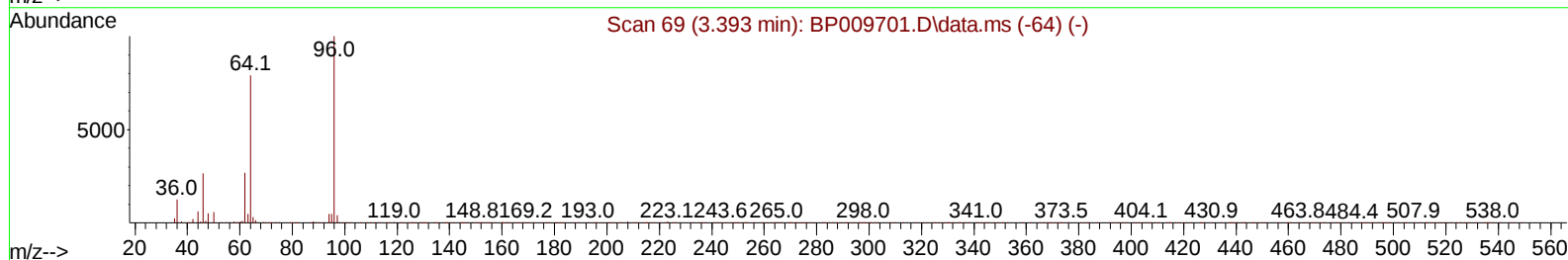
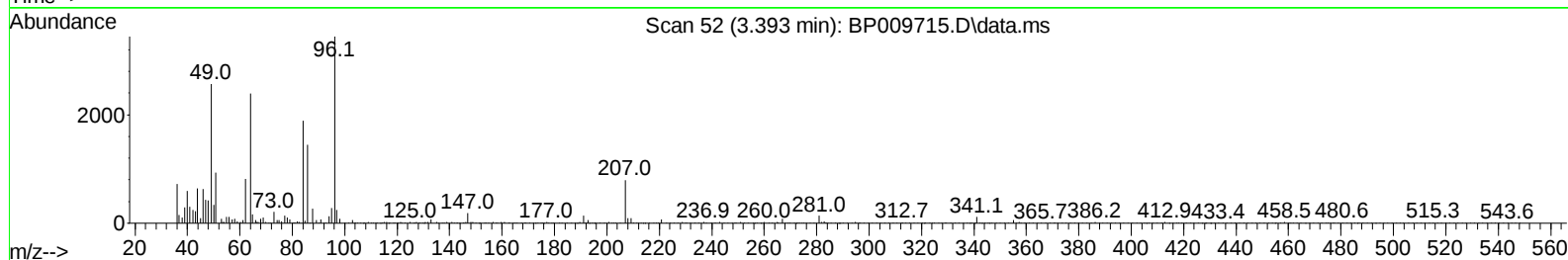
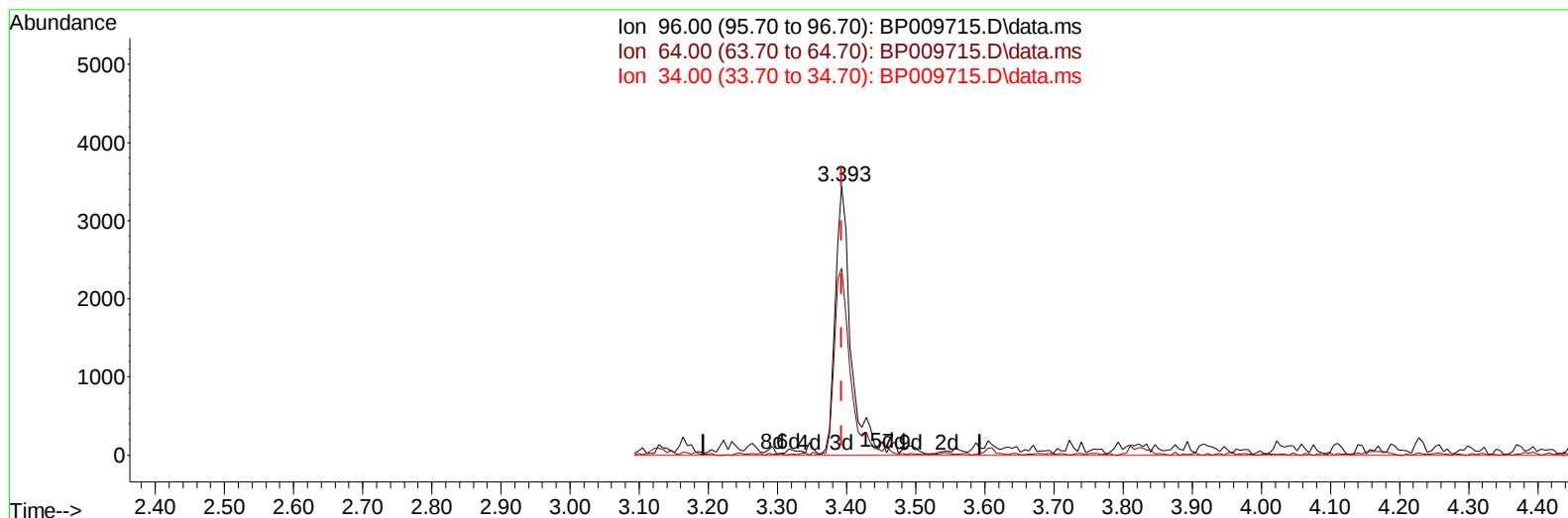
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TIC: BP009715.D\data.ms

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

3.393min (-0.000) 2.56 ng/uL m

response 5293

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	69.49#
34.00	0.00	0.00
0.00	0.00	0.00

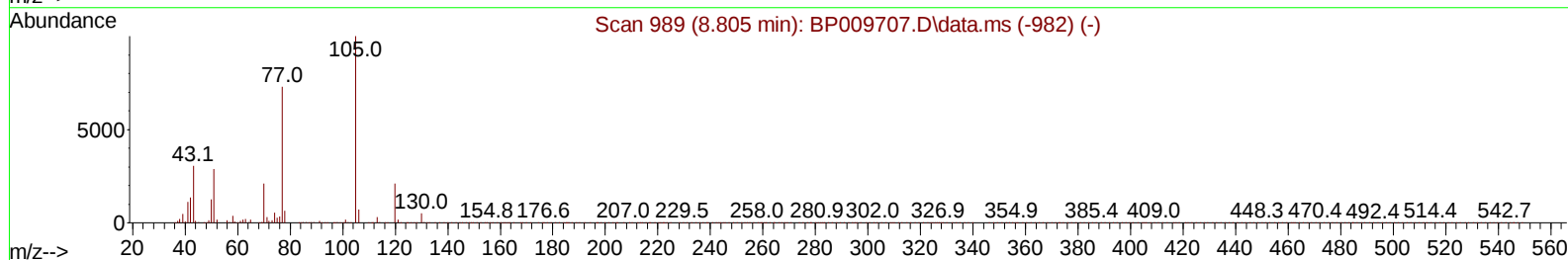
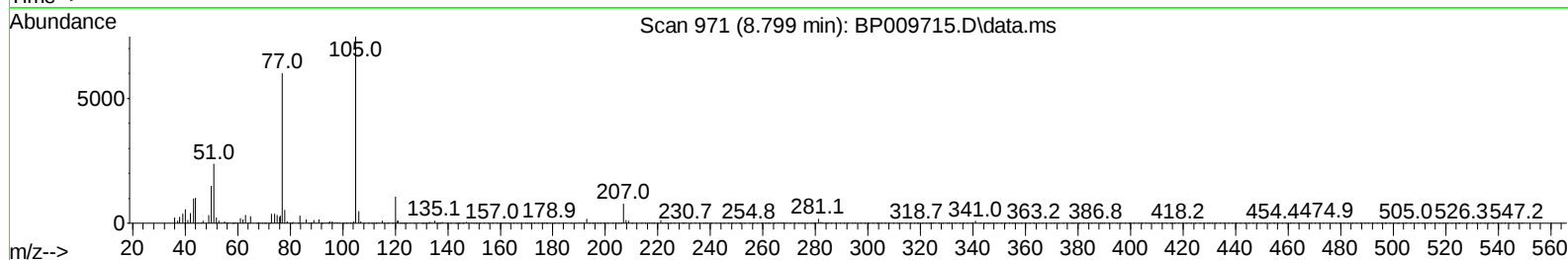
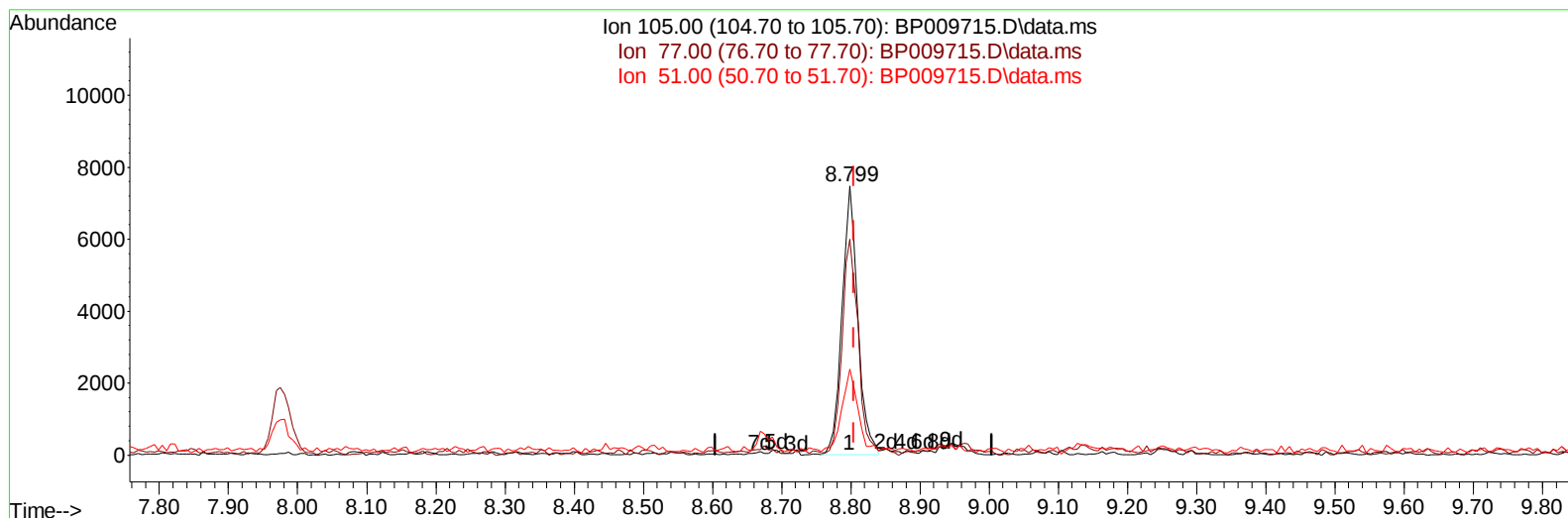
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TIC: BP009715.D\data.ms

(16) Acetophenone

8.799min (-0.006) 1.11 ng/u1

response 11846

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	66.80	80.39#
51.00	17.30	31.85#
0.00	0.00	0.00

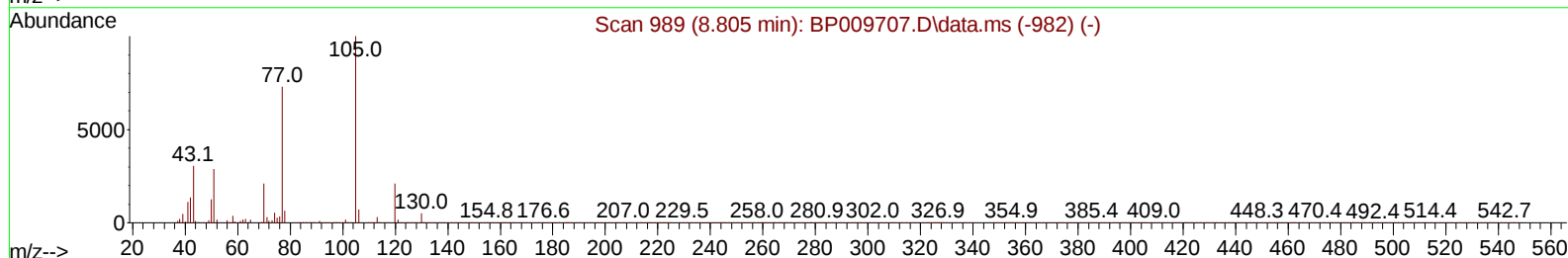
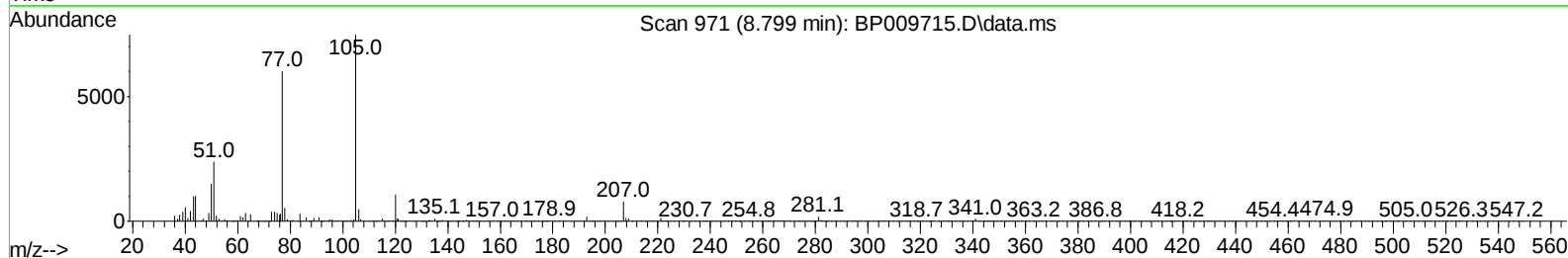
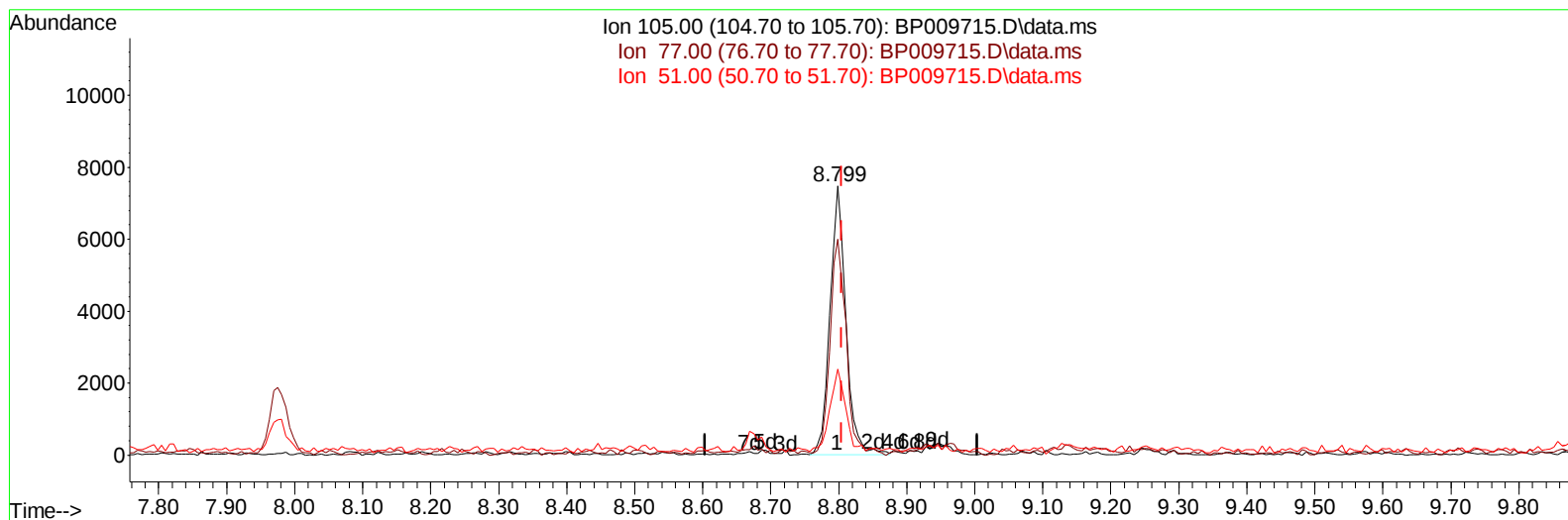
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TIC: BP009715.D\data.ms

(16) Acetophenone

8.799min (-0.006) 1.14 ng/ul m

response 12114

Ion	Exp%	Act%
105.00	100.00	100.00
77.00	66.80	80.39#
51.00	17.30	31.85#
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	92110	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	401889	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.610	164	303471	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.357	188	699725	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.439	240	771969	20.000	ng/ul	# 0.00
88) Perylene-d12	23.915	264	786822	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	5293m	2.565	ng/uL	0.00
4) Pyridine-d5	3.811	84	41417	7.278	ng/ul	0.00
7) Phenol-d5	7.122	99	88998	11.005	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	61155	12.703	ng/ul	0.00
11) 2-Chlorophenol-d4	7.505	132	71095	11.655	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	82145	12.168	ng/ul	0.00
21) Nitrobenzene-d5	9.134	128	37007	12.769	ng/ul	0.00
24) 2-Nitrophenol-d4	9.857	143	39327	12.377	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.398	165	75809	11.402	ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131	108356	11.442	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	336474	14.115	ng/ul	0.00
49) Acenaphthylene-d8	14.298	160	352281	12.474	ng/ul	0.00
54) 4-Nitrophenol-d4	14.781	143	48026	11.109	ng/ul	0.00
60) Fluorene-d10	15.598	176	282348	14.080	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	39226	9.463	ng/ul	0.00
73) Anthracene-d10	17.457	188	528261	15.768	ng/ul	0.00
81) Pyrene-d10	19.686	212	719218	17.059	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.757	264	726867	17.739	ng/ul	0.00
Target Compounds						
16) Acetophenone	8.799	105	12114m	1.139	ng/ul	Qvalue

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

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