

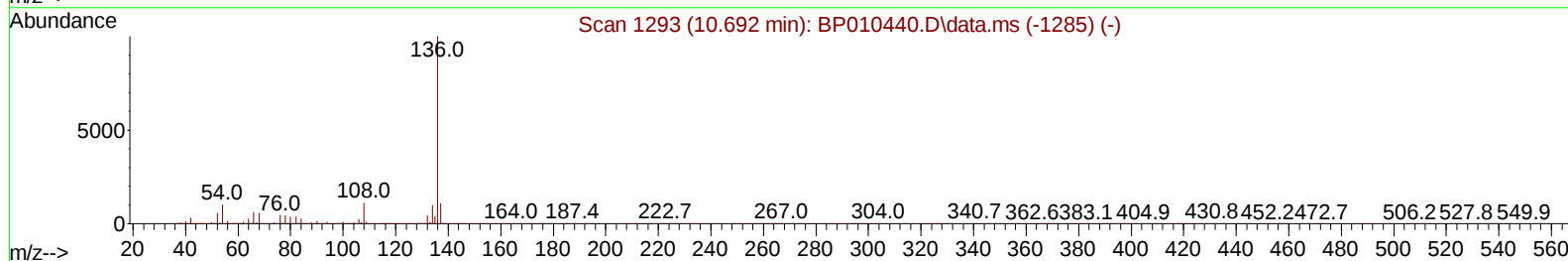
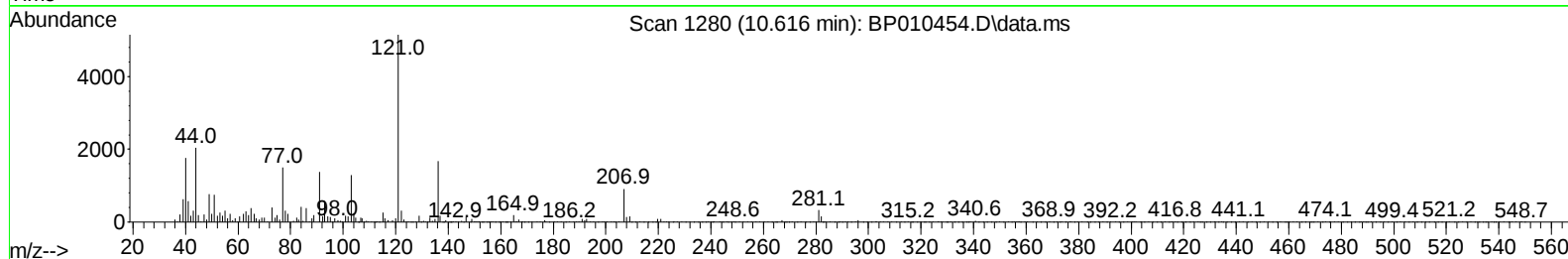
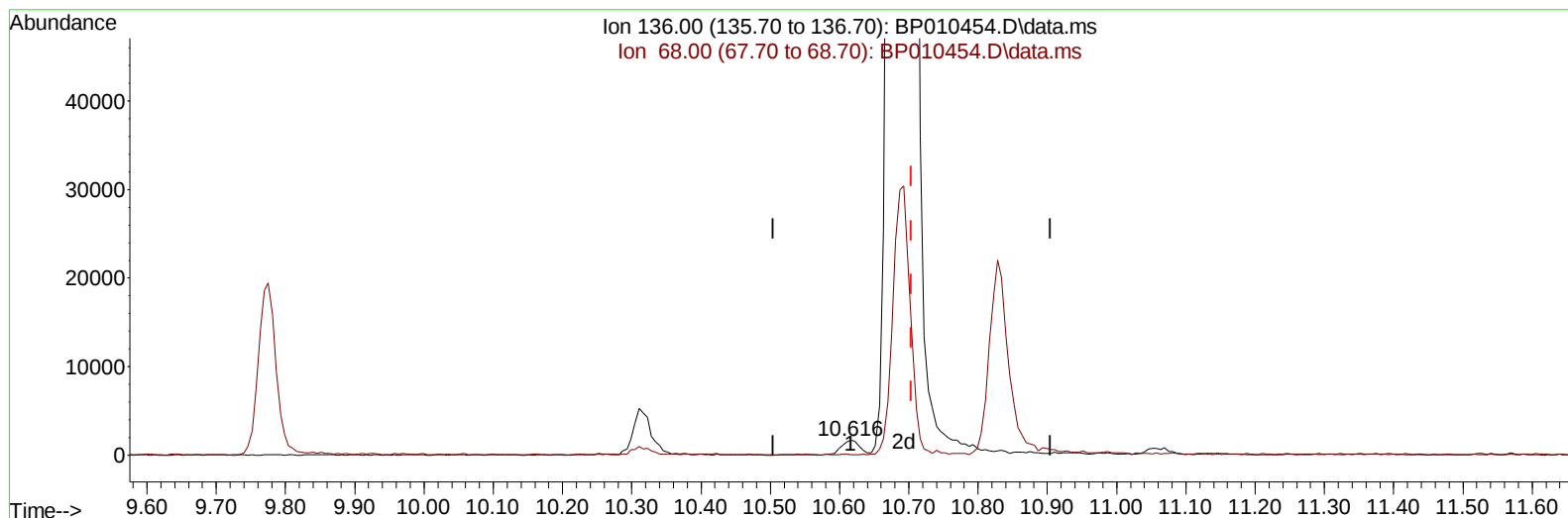
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052022\
 Data File : BP010454.D
 Acq On : 21 May 2022 01:56
 Operator : CG/JU
 Sample : N2843-08
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW4H8

Manual IntegrationsAPPROVED

Quant Time: May 21 03:50:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 13 15:04:31 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/23/2022
 Supervised By :mohammad ahmed 05/25/2022



TIC: BP010454.D\data.ms

(20) Naphthalene-d8 (I)

10.616min (-0.088) 20.00 ng/u1

response 3094

Ion	Exp%	Act%
136.00	100.00	100.00
68.00	5.10	4.30
0.00	0.00	0.00
0.00	0.00	0.00

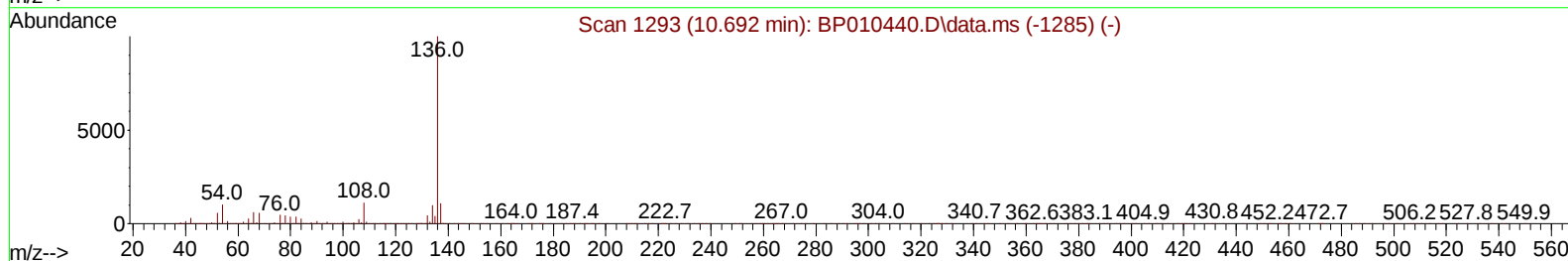
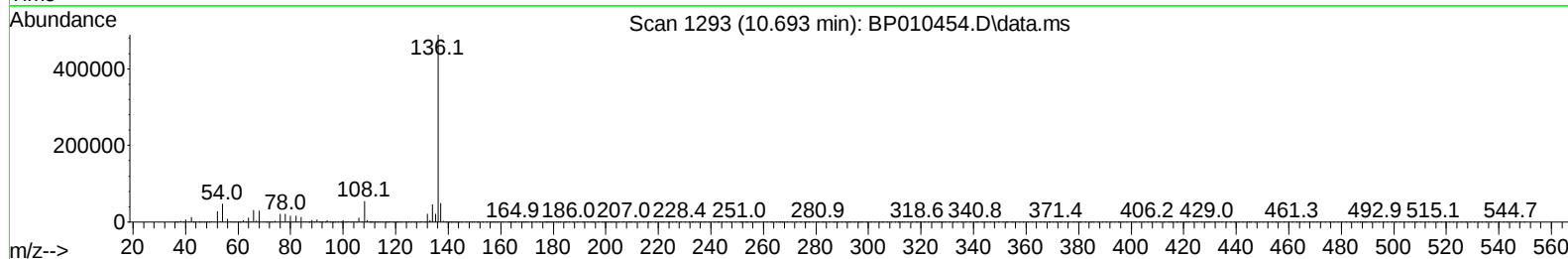
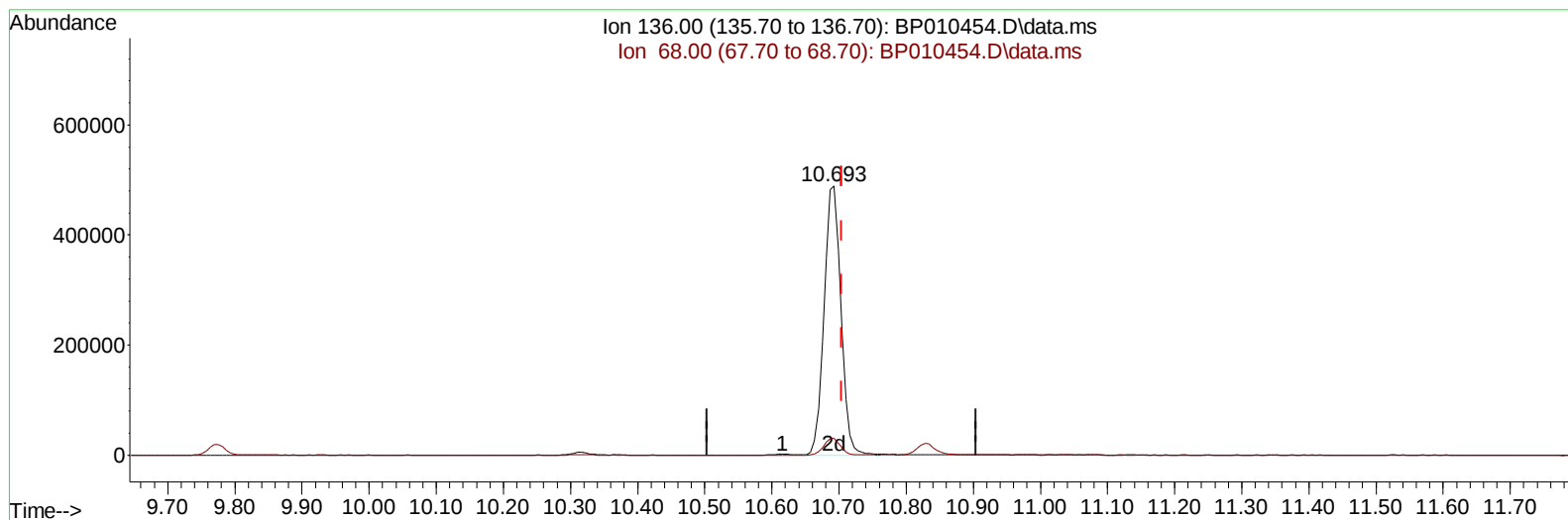
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(20) Naphthalene-d8 (I)

10.693min (-0.011) 20.00 ng/ul m

response 856158

Ion	Exp%	Act%
136.00	100.00	100.00
68.00	5.10	6.22#
0.00	0.00	0.00
0.00	0.00	0.00

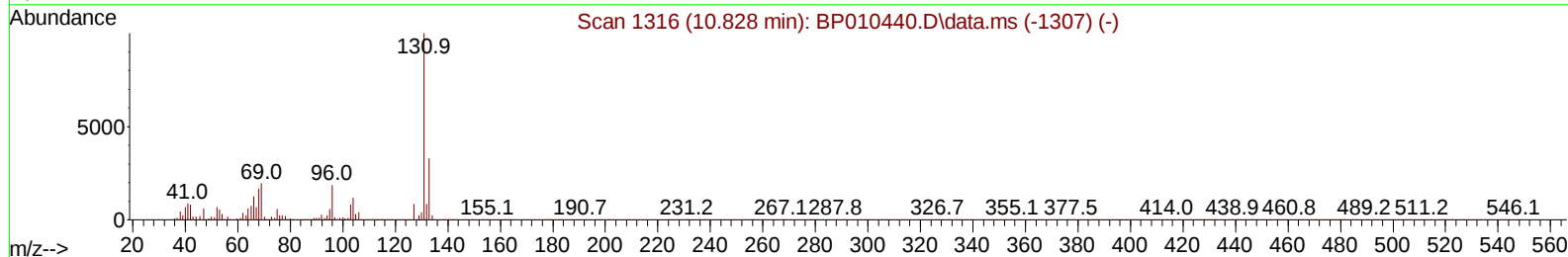
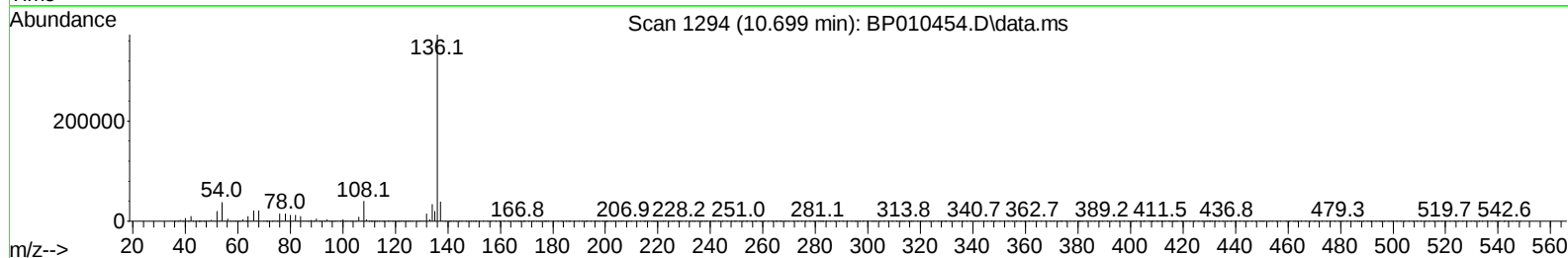
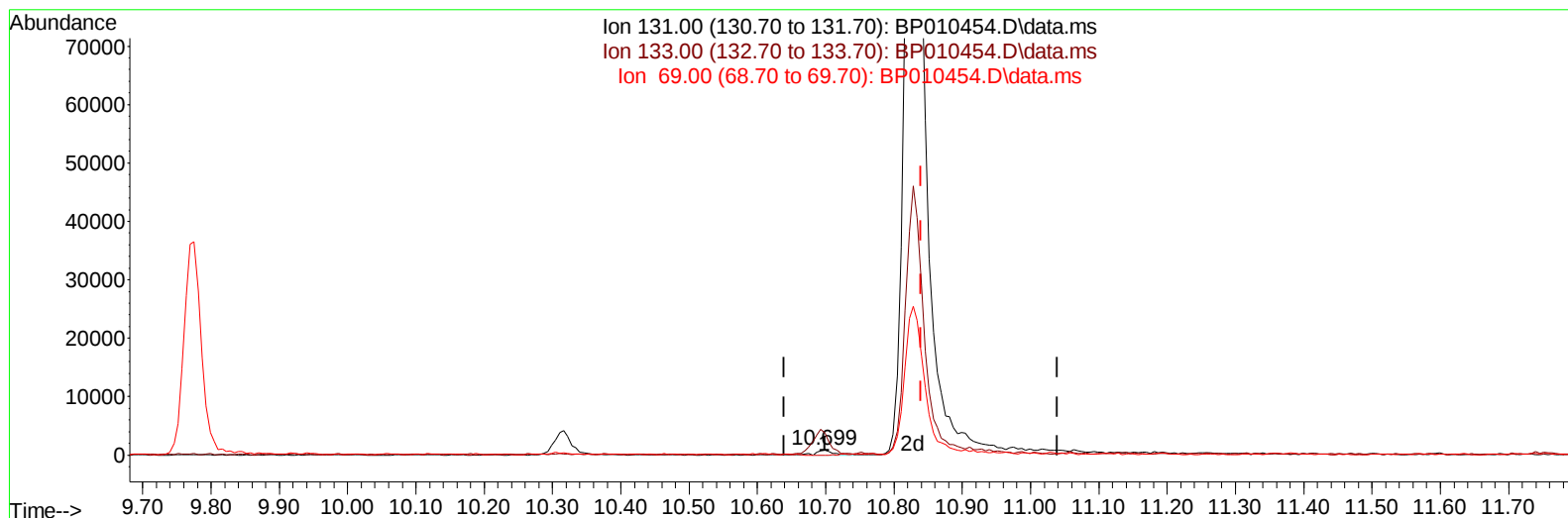
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TIC: BP010454.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.699min (-0.141) 0.06 ng/u1

response 1122

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	454.13#
69.00	26.10	0.00#
0.00	0.00	0.00

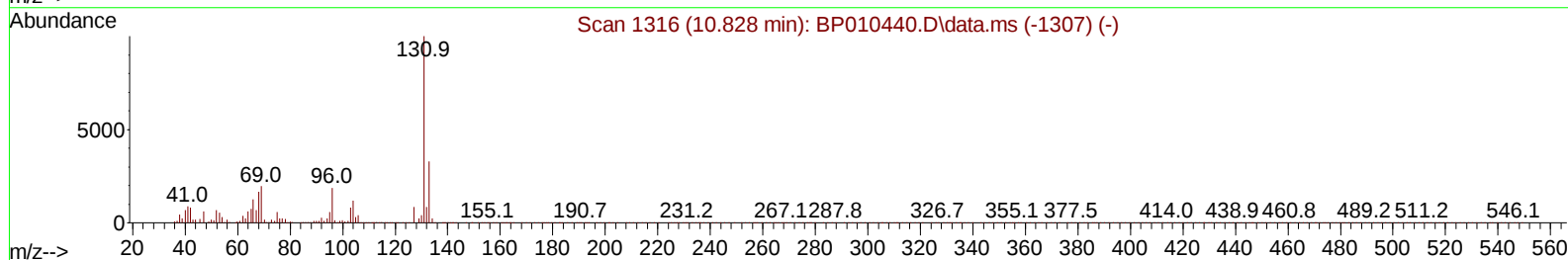
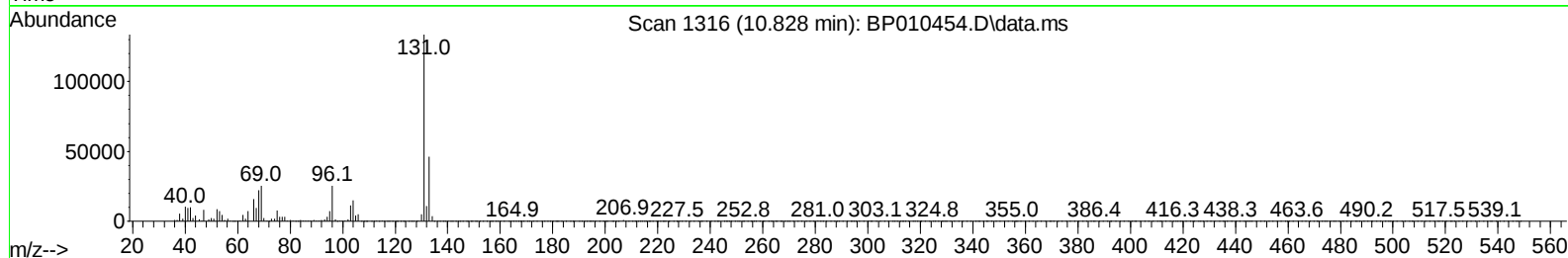
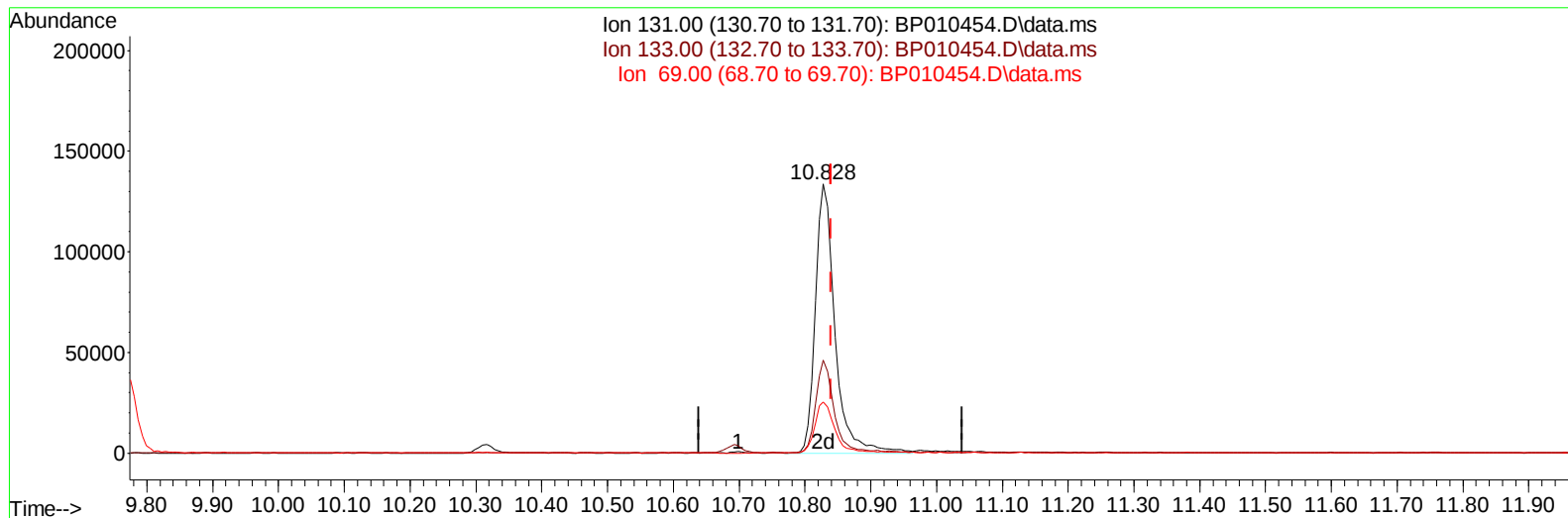
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Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP010454.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.828min (-0.011) 13.87 ng/ul m

response 273783

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	34.45
69.00	26.10	19.01#
0.00	0.00	0.00

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 ALS Vial : 29 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	189172	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	856158m	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.522	164	587206	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.275	188	1212912	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.357	240	1076610	20.000	ng/ul	0.00
88) Perylene-d12	23.780	264	991726	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	10887	2.453	ng/uL	0.00
4) Pyridine-d5	3.764	84	81772	6.489	ng/ul	0.00
7) Phenol-d5	7.058	99	244184	15.379	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.217	67	173492	16.263	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.422	132	196166	16.146	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	210258	15.582	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	108917	16.438	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.775	143	114893	16.362	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	198698	14.965	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.828	131	273783m	13.868	ng/ul	-0.01
46) Dimethylphthalate-d6	13.934	166	722216	16.693	ng/ul	0.00
49) Acenaphthylene-d8	14.210	160	851715	17.067	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.722	143	86098	11.181	ng/ul	0.00
60) Fluorene-d10	15.516	176	655789	17.423	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	81725	11.521	ng/ul	-0.01
73) Anthracene-d10	17.375	188	1027430	18.707	ng/ul	0.00
81) Pyrene-d10	19.604	212	1219720	21.333	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.622	264	959960	19.277	ng/ul	-0.01
Target Compounds						
					Qvalue	
16) Acetophenone	8.716	105	32336	1.525	ng/ul#	78

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

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