

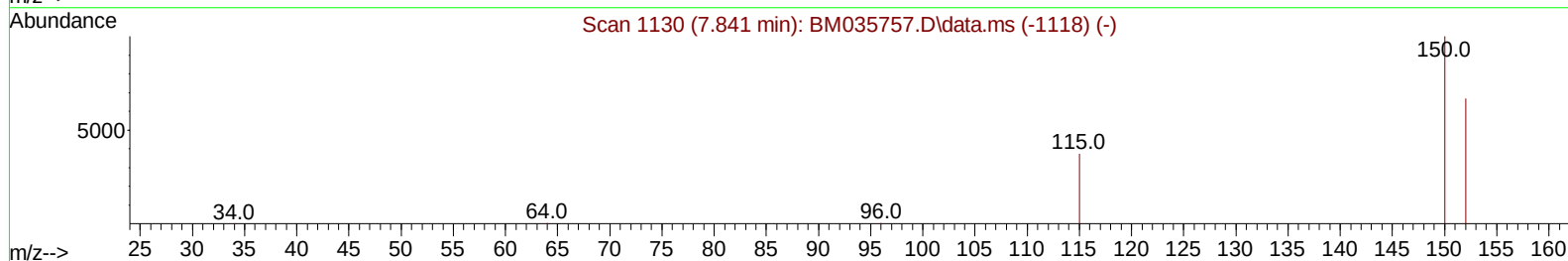
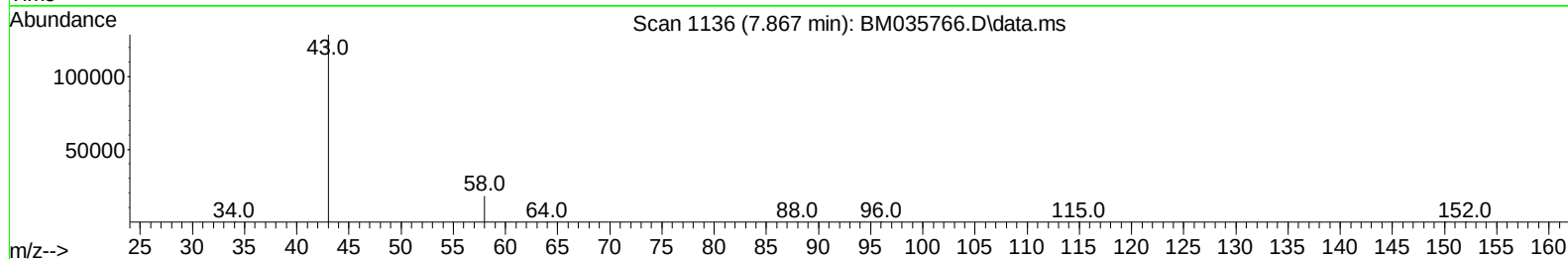
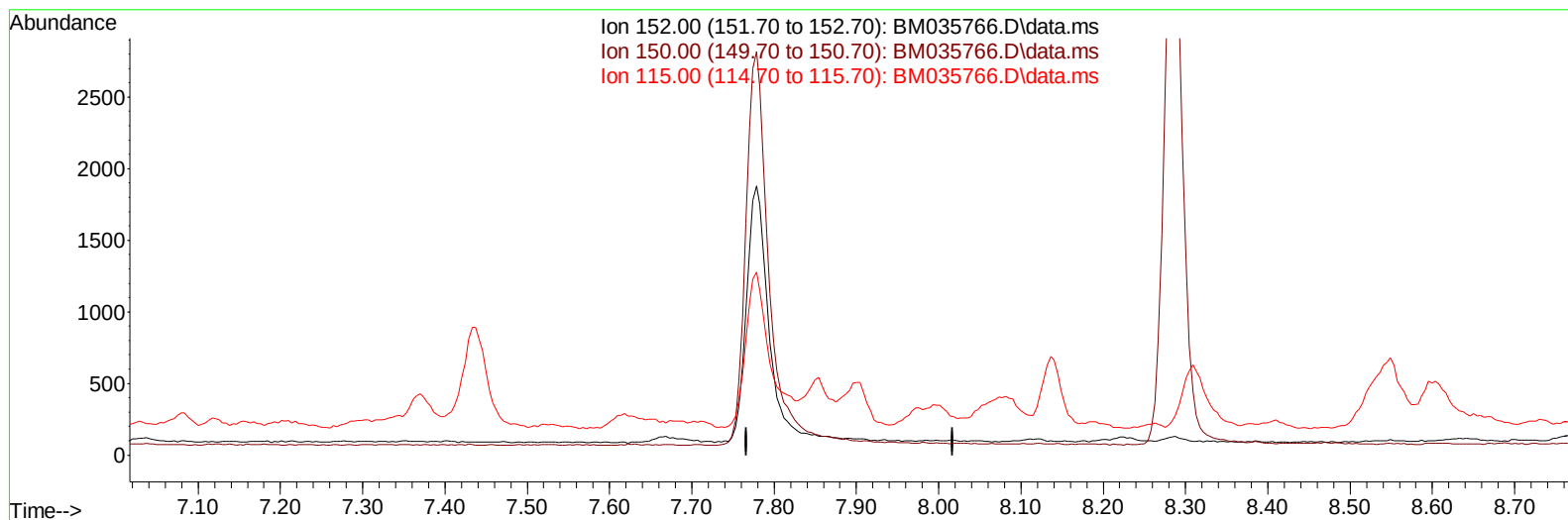
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062822\
Data File : BM035766.D
Acq On : 29 Jun 2022 18:15
Operator : CG/JU
Sample : N3374-02MS
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4S6MSA

Manual IntegrationsAPPROVED

Quant Time: Jun 30 03:57:51 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 29 17:48:53 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/30/2022
Supervised By :mohammad ahmed 07/01/2022



TIC: BM035766.D\data.ms

(1) 1,4-Dichlorobenzene-d4 (I)

7.867min (-7.867) 0.00 ng/u1

response 0

Ion	Exp%	Act%
152.00	100.00	0.00
150.00	147.40	0.00#
115.00	74.40	0.00#
0.00	0.00	0.00

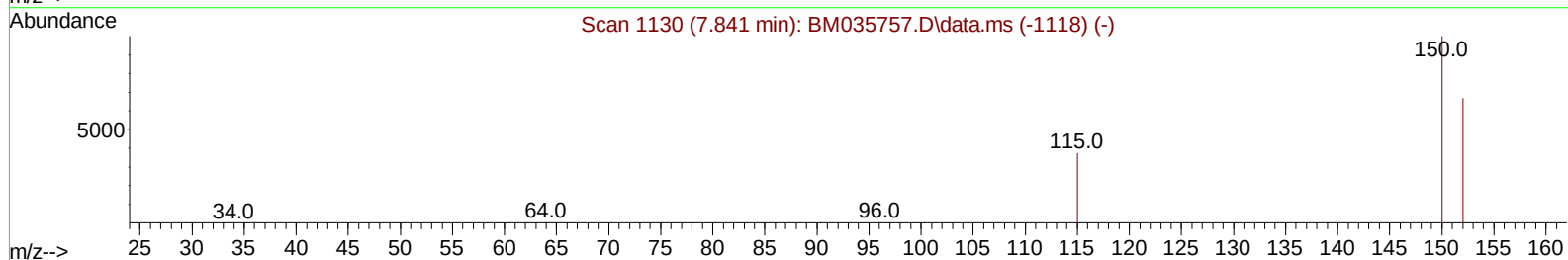
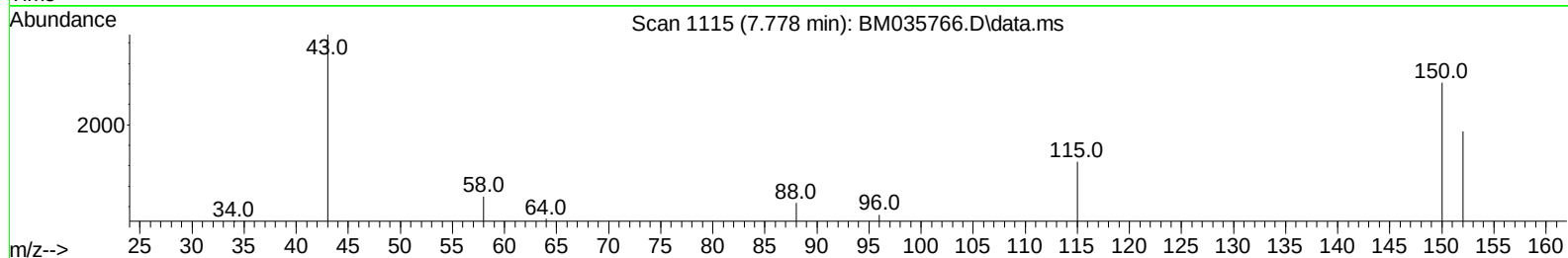
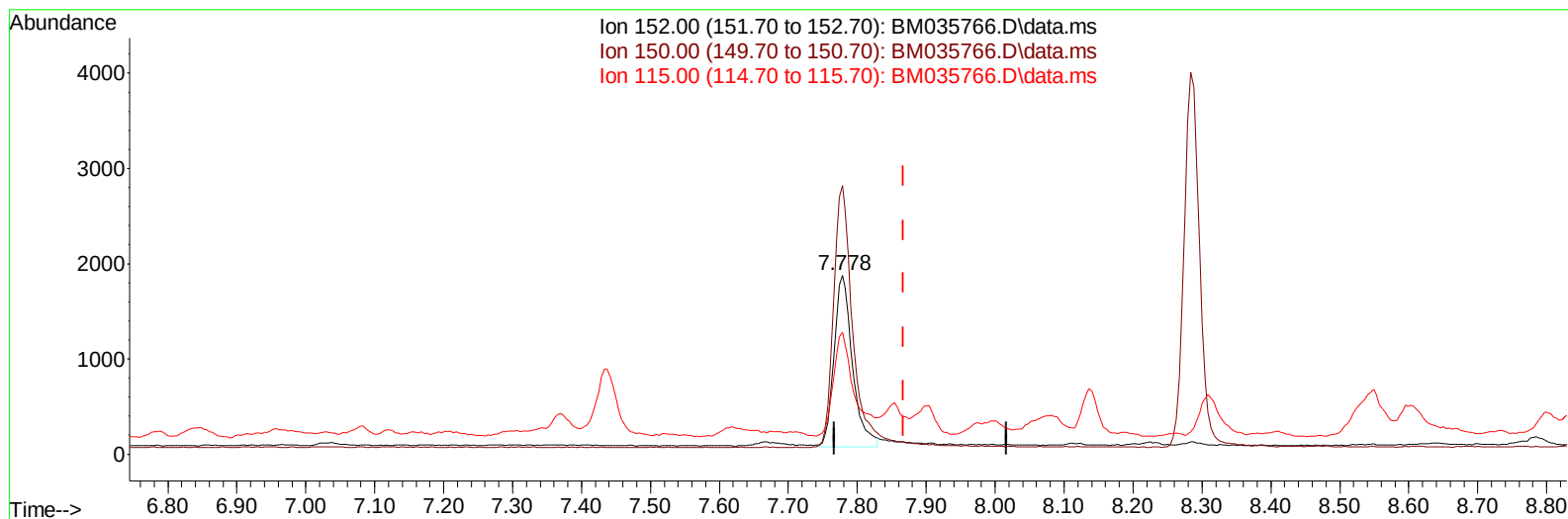
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(1) 1,4-Dichlorobenzene-d4 (I)

7.778min (-0.089) 0.40 ng/ul m

response 3366

Ion	Exp%	Act%
152.00	100.00	100.00
150.00	147.40	150.05
115.00	74.40	67.94
0.00	0.00	0.00

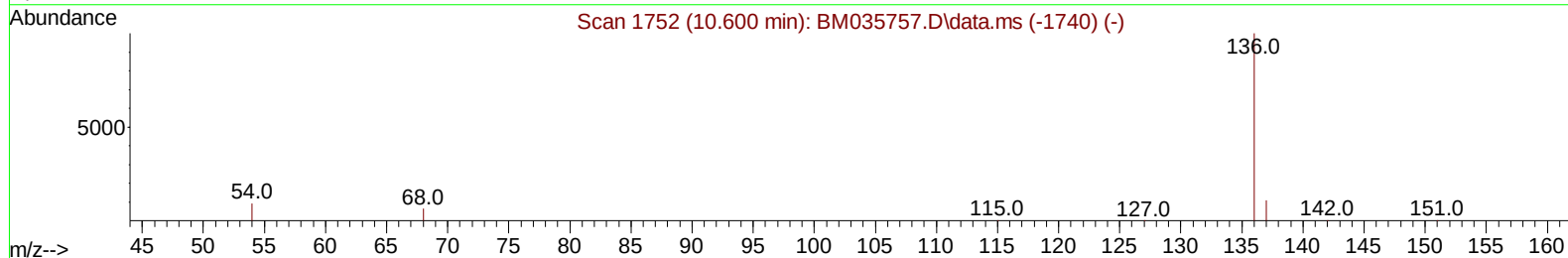
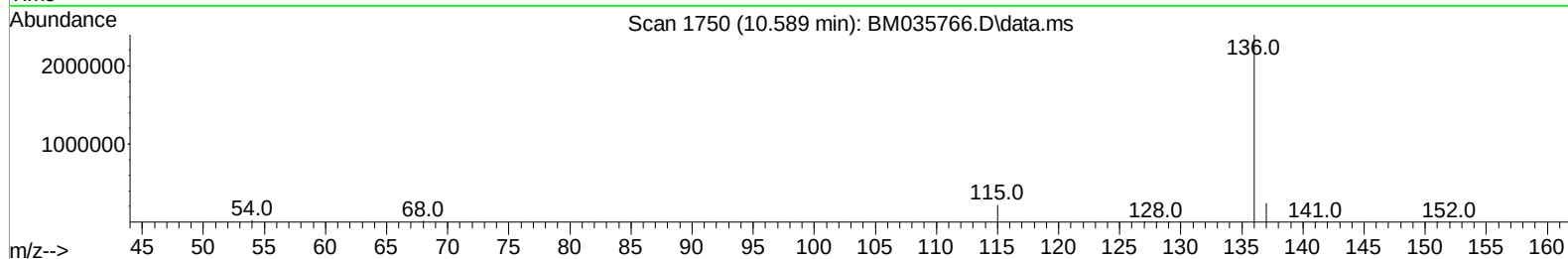
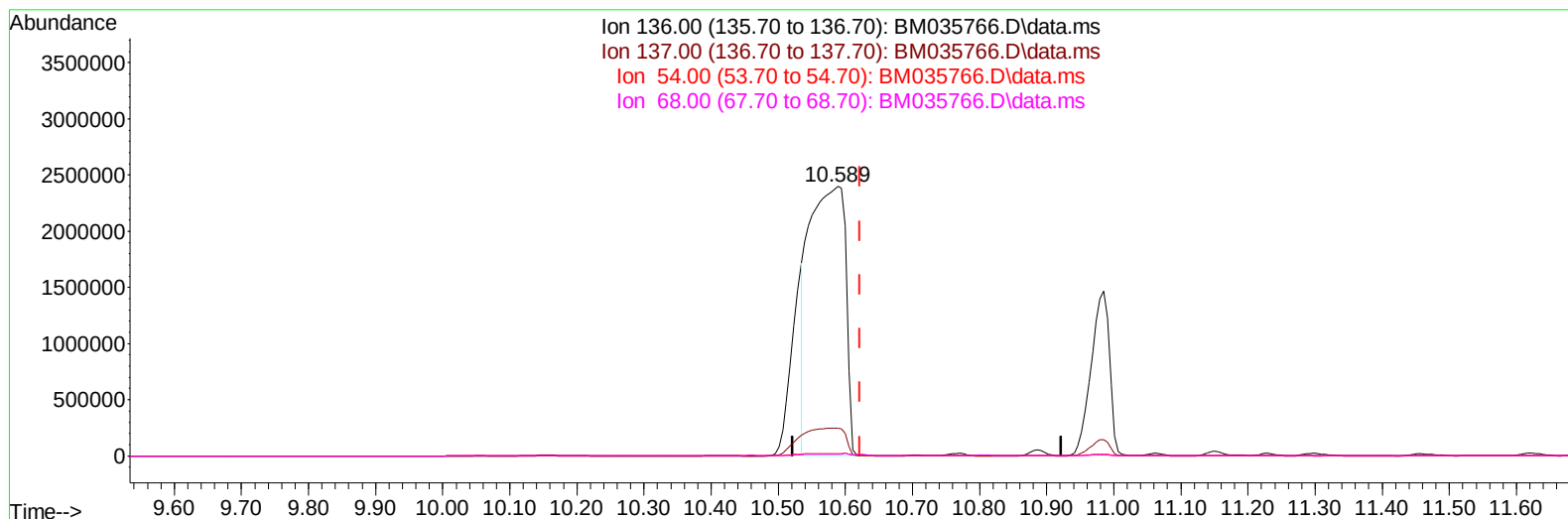
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TIC: BM035766.D\data.ms

(4) Naphthalene-d8 (I)

10.589min (-0.033) 0.40 ng/u1

response 9086004

Ion	Exp%	Act%
136.00	100.00	100.00
137.00	12.20	10.27
54.00	12.20	0.89#
68.00	8.00	0.57#

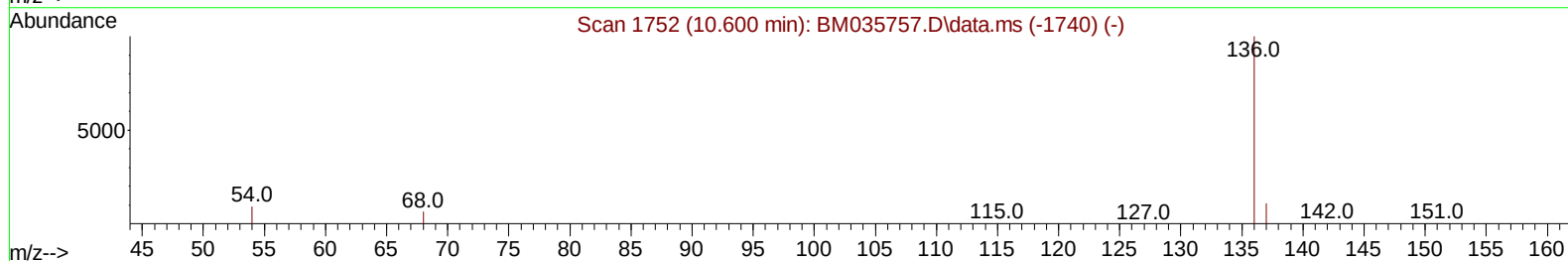
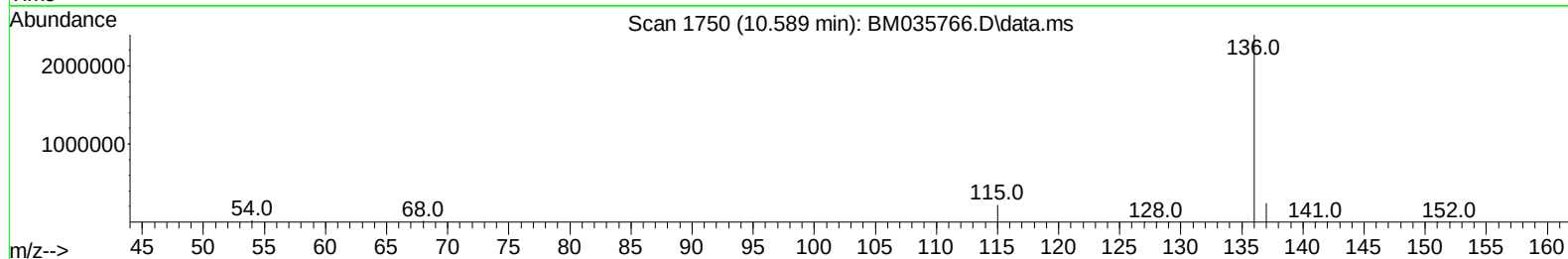
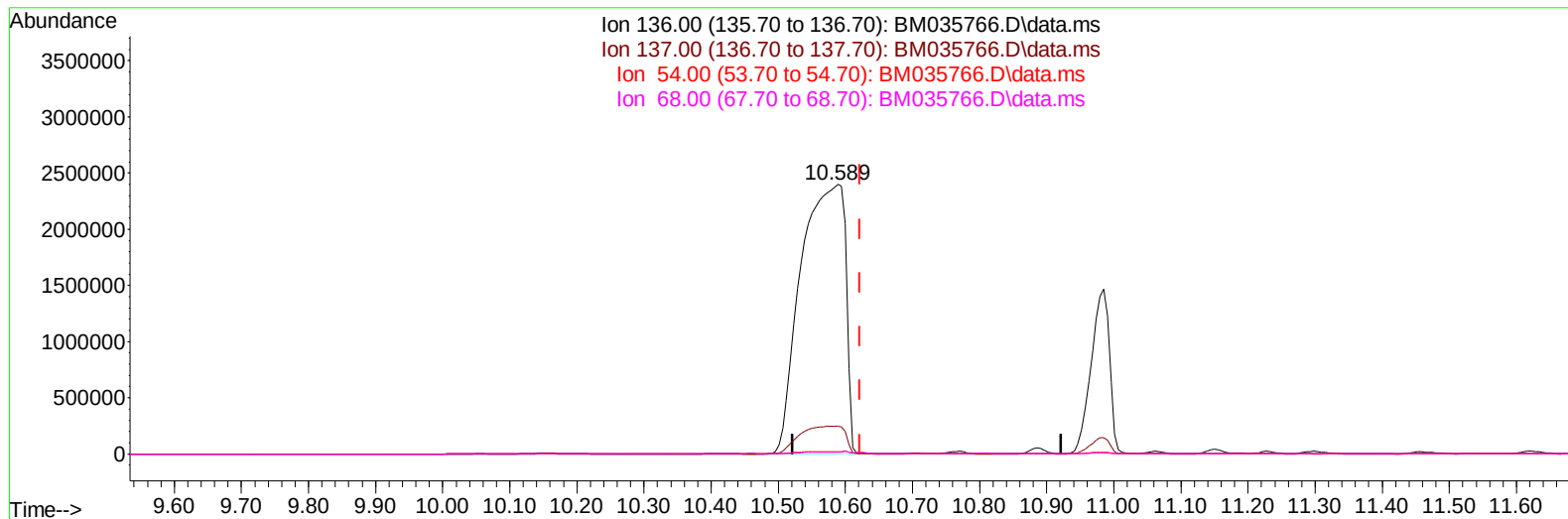
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Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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TIC: BM035766.D\data.ms

(4) Naphthalene-d8 (I)

10.589min (-0.033) 0.40 ng/ul m

response 11074222

Ion	Exp%	Act%
136.00	100.00	100.00
137.00	12.20	10.27
54.00	12.20	0.89#
68.00	8.00	0.57#

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.778	152		3366m	0.400	ng/ul	-0.09
4) Naphthalene-d8	10.589	136	11074222m		0.400	ng/ul	-0.03
9) Acenaphthene-d10	14.418	164		7282	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.165	188		11817	0.400	ng/ul	#-0.03
17) Chrysene-d12	21.357	240		11621	0.400	ng/ul	#-0.02
23) Perylene-d12	23.677	264		12279	0.400	ng/ul	-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.217	96		1250	0.274	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.167	152		11284	0.001	ng/ul	-0.06
18) Fluoranthene-d10	19.196	212		10851	0.277	ng/ul	-0.02
Target Compounds					Qvalue		
2) 1,4-Dioxane	3.247	88		19471	4.310	ng/ul#	1
11) Acenaphthene	14.478	153		31705	1.078	ng/ul	98
12) Fluorene	15.468	166		12199	0.360	ng/ul#	95
14) Pentachlorophenol	16.845	266		999	0.535	ng/ul	90
15) Phenanthrene	17.203	178		21949	0.513	ng/ul#	89
16) Anthracene	17.296	178		24111	0.678	ng/ul#	89
19) Fluoranthene	19.224	202		20322	0.349	ng/ul#	86
20) Pyrene	19.591	202		19596	0.300	ng/ul#	90
21) Benzo(a)anthracene	21.342	228		14780	0.368	ng/ul#	80
22) Chrysene	21.395	228		15871	0.307	ng/ul#	80
24) Benzo(b)fluoranthene	22.976	252		16603	0.315	ng/ul	64
25) Benzo(k)fluoranthene	23.022	252		15924	0.296	ng/ul#	65
26) Benzo(a)pyrene	23.578	252		16008	0.298	ng/ul#	73
27) Indeno(1,2,3-cd)pyrene	26.085	276		19360	0.330	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.088	278		15773	0.340	ng/ul#	70
29) Benzo(g,h,i)perylene	26.826	276		17397	0.322	ng/ul#	79

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

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