

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102218\
 Data File : BM017279.D
 Acq On : 22 Oct 2018 15:05
 Operator : MJ/SJ
 Sample : J5585-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 23 07:45:02 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM101718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 17 14:23:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	356427	20.00	ng	-0.02
21) Naphthalene-d8	10.43	136	1485623	20.00	ng	-0.02
39) Acenaphthene-d10	14.30	164	833821	20.00	ng	-0.02
64) Phenanthrene-d10	17.05	188	1715852	20.00	ng	-0.02
76) Chrysene-d12	21.24	240	1509396	20.00	ng	-0.02
87) Perylene-d12	23.46	264	1437974	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.28	112	2271853	107.81	ng	-0.01
7) Phenol-d6	6.85	99	3204748	111.66	ng	-0.01
23) Nitrobenzene-d5	8.82	82	1964161	77.32	ng	-0.02
42) 2,4,6-Tribromophenol	15.80	330	1189436	105.48	ng	-0.02
45) 2-Fluorobiphenyl	12.92	172	4701683	75.03	ng	-0.02
79) Terphenyl-d14	19.69	244	5747869	85.83	ng	-0.01
Target Compounds						
10) Phenol	6.87	94	280451	9.079	ng	98
50) Dimethylphthalate	13.76	163	578637	8.995	ng	99

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

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