

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM102318\
 Data File : BM017307.D
 Acq On : 24 Oct 2018 08:32
 Operator : MJ/SJ
 Sample : J5623-12
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 101918-DBRI-E-U-B

Quant Time: Oct 24 09:19:48 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	263456	20.00	ng	-0.02
21) Naphthalene-d8	10.43	136	1110120	20.00	ng	-0.02
39) Acenaphthene-d10	14.30	164	643801	20.00	ng	-0.02
64) Phenanthrene-d10	17.05	188	1514011	20.00	ng	-0.02
76) Chrysene-d12	21.24	240	1702308	20.00	ng	-0.02
87) Perylene-d12	23.46	264	1481687	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	680873	43.71	ng	-0.01
7) Phenol-d6	6.84	99	583000	27.48	ng	-0.02
23) Nitrobenzene-d5	8.81	82	2031327	107.02	ng	-0.02
42) 2,4,6-Tribromophenol	15.80	330	911594	104.70	ng	-0.02
45) 2-Fluorobiphenyl	12.92	172	4932558	101.94	ng	-0.02
79) Terphenyl-d14	19.69	244	6838716	90.55	ng	-0.01
Target Compounds						
10) Phenol	6.86	94	75202	3.294	ng	95
50) Dimethylphthalate	13.76	163	107457	2.163	ng	99

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

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