

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM102318\
 Data File : BM017298.D
 Acq On : 23 Oct 2018 23:18
 Operator : MJ/SJ
 Sample : J5604-08
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MS-VB-25868-BOX-1

Quant Time: Oct 24 07:56:12 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	307308	20.00	ng	-0.01
21) Naphthalene-d8	10.43	136	1113588	20.00	ng	-0.02
39) Acenaphthene-d10	14.30	164	579203	20.00	ng	-0.02
64) Phenanthrene-d10	17.05	188	1188909	20.00	ng	-0.02
76) Chrysene-d12	21.24	240	1295460	20.00	ng	-0.02
87) Perylene-d12	23.46	264	1405025	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	62594	3.44	ng	0.00
7) Phenol-d6	6.85	99	677927	27.40	ng	-0.01
23) Nitrobenzene-d5	8.82	82	1532472	80.49	ng	-0.02
42) 2,4,6-Tribromophenol	15.80	330	896956	114.51	ng	-0.02
45) 2-Fluorobiphenyl	12.92	172	3390041	77.88	ng	-0.02
79) Terphenyl-d14	19.69	244	4133407	71.92	ng	-0.01
Target Compounds						Qvalue
9) Benzaldehyde	6.82	77	64958	4.117	ng	96
10) Phenol	6.87	94	127408	4.784	ng	99
27) 2,4-Dimethylphenol	9.66	122	129275	9.299	ng	# 88
32) Benzoic acid	9.73	122	35508	3.423	ng	# 85
50) Dimethylphthalate	13.76	163	251660	5.632	ng	99

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

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