

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070618\
 Data File : BM015810.D
 Acq On : 06 Jul 2018 20:07
 Operator : SJ/JU
 Sample : J3864-02RE
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EPE-108-6A(19-21)RE

Quant Time: Jul 07 01:31:37 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 05 17:39:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.30	152	112592	20.00	ng	-0.01
21) Naphthalene-d8	11.13	136	440792	20.00	ng	-0.01
38) Acenaphthene-d10	14.93	164	236618	20.00	ng	0.00
63) Phenanthrene-d10	17.65	188	470568	20.00	ng	0.00
75) Chrysene-d12	21.76	240	335611	20.00	ng	-0.01
86) Perylene-d12	24.16	264	318485	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.81	112	936673	142.67	ng	0.00
7) Phenol-d6	7.46	99	1226312	136.44	ng	0.00
23) Nitrobenzene-d5	9.49	82	812706	89.96	ng	-0.01
41) 2,4,6-Tribromophenol	16.40	330	400406	129.51	ng	-0.01
44) 2-Fluorobiphenyl	13.55	172	1473019	77.29	ng	-0.01
78) Terphenyl-d14	20.22	244	1464331	80.93	ng	0.00
Target Compounds						
10) Phenol	7.49	94	21886	2.423	ng	88
49) Dimethylphthalate	14.37	163	309624	14.837	ng	99

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

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