

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011719\
 Data File : BM018622.D
 Acq On : 17 Jan 2019 19:58
 Operator : JU/SJ
 Sample : K1153-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RBR200044-RBR200027

Quant Time: Jan 18 06:03:28 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 10 05:57:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	391758	20.00	ng	-0.02
21) Naphthalene-d8	10.53	136	1596636	20.00	ng	-0.02
39) Acenaphthene-d10	14.39	164	908363	20.00	ng	-0.02
64) Phenanthrene-d10	17.14	188	1954486	20.00	ng	-0.02
76) Chrysene-d12	21.33	240	1387812	20.00	ng	-0.02
87) Perylene-d12	23.60	264	1293857	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	2343456	110.46	ng	-0.02
7) Phenol-d6	6.92	99	3056945	113.44	ng	-0.02
23) Nitrobenzene-d5	8.90	82	1861609	67.59	ng	-0.02
42) 2,4,6-Tribromophenol	15.89	330	1308841	113.16	ng	-0.02
45) 2-Fluorobiphenyl	13.00	172	4154485	59.68	ng	-0.02
79) Terphenyl-d14	19.77	244	4869848	73.97	ng	-0.02
Target Compounds						
10) Phenol	6.94	94	56048	2.047	ng	81
50) Dimethylphthalate	13.84	163	352102	4.720	ng	99
71) Phenanthrene	17.18	178	279908	2.692	ng	99
75) Fluoranthene	19.20	202	652134	5.441	ng	98
78) Pyrene	19.56	202	558332	6.787	ng	100
81) Benzo(a)anthracene	21.31	228	237071	2.900	ng	98
83) Chrysene	21.37	228	244619	3.098	ng	97
84) Bis(2-ethylhexyl)phthalate	21.24	149	204017	4.025	ng	99
88) Benzo(b)fluoranthene	22.91	252	297860	4.055	ng	98
90) Benzo(a)pyrene	23.50	252	198569	2.824	ng	99

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

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