

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060519\
 Data File : BM020755.D
 Acq On : 06 Jun 2019 02:30
 Operator : HP/JU
 Sample : K3183-09
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-3

Quant Time: Jun 06 04:10:27 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM053119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 31 15:48:49 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	293168	20.00	ng	-0.01
21) Naphthalene-d8	10.61	136	1141284	20.00	ng	-0.02
39) Acenaphthene-d10	14.46	164	592092	20.00	ng	-0.02
64) Phenanthrene-d10	17.20	188	1227975	20.00	ng	-0.02
76) Chrysene-d12	21.37	240	1185487	20.00	ng	-0.02
87) Perylene-d12	23.66	264	1439562	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	1717761	103.07	ng	-0.01
7) Phenol-d6	6.98	99	2339258	95.32	ng	-0.01
23) Nitrobenzene-d5	8.98	82	1428098	64.83	ng	-0.02
42) 2,4,6-Tribromophenol	15.95	330	742128	101.16	ng	-0.02
45) 2-Fluorobiphenyl	13.08	172	2944281	66.11	ng	-0.01
79) Terphenyl-d14	19.82	244	3514510	49.59	ng	-0.02
Target Compounds				Qvalue		
50) Dimethylphthalate	13.92	163	479469	9.478	ng	100
71) Phenanthrene	17.25	178	285584	4.103	ng	99
75) Fluoranthene	19.26	202	464934	6.006	ng	99
78) Pyrene	19.62	202	409560	4.502	ng	99
81) Benzo(a)anthracene	21.36	228	251337	3.044	ng	99
83) Chrysene	21.41	228	206261	2.628	ng	99
88) Benzo(b)fluoranthene	22.97	252	269627	2.834	ng	95
90) Benzo(a)pyrene	23.56	252	191949	2.224	ng	99

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

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