

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022619\
 Data File : BM018931.D
 Acq On : 26 Feb 2019 16:09
 Operator : JU/SJ
 Sample : K1597-15
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-23

Quant Time: Feb 27 00:36:36 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 11 16:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	254439	20.00	ng	-0.02
21) Naphthalene-d8	10.48	136	920259	20.00	ng	-0.02
39) Acenaphthene-d10	14.34	164	439308	20.00	ng	-0.02
64) Phenanthrene-d10	17.10	188	855729	20.00	ng	-0.02
76) Chrysene-d12	21.29	240	956131	20.00	ng	-0.02
87) Perylene-d12	23.54	264	1110224	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	2297354	147.94	ng	-0.02
7) Phenol-d6	6.87	99	3064283	140.07	ng	-0.01
23) Nitrobenzene-d5	8.86	82	1985082	99.74	ng	-0.02
42) 2,4,6-Tribromophenol	15.84	330	772834	122.04	ng	-0.02
45) 2-Fluorobiphenyl	12.96	172	2625672	79.34	ng	-0.02
79) Terphenyl-d14	19.73	244	2996318	64.65	ng	-0.01
Target Compounds						
10) Phenol	6.90	94	196805	8.550	ng	Qvalue 98
50) Dimethylphthalate	13.80	163	200798	5.485	ng	98
71) Phenanthrene	17.14	178	111694	2.429	ng	99
75) Fluoranthene	19.16	202	210493	3.962	ng	99
78) Pyrene	19.53	202	210269	3.787	ng	99
81) Benzo(a)anthracene	21.27	228	111759	2.005	ng	97
83) Chrysene	21.33	228	111571	2.089	ng	98
88) Benzo(b)fluoranthene	22.87	252	156802	2.469	ng	# 92

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

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