

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090518\
 Data File : BM016671.D
 Acq On : 05 Sep 2018 14:53
 Operator : SJ/JU
 Sample : J4764-01RE
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SPRE

Quant Time: Sep 05 17:13:00 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 05 16:42:07 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.97	152	60822	20.00	ng	0.00
21) Naphthalene-d8	10.77	136	209245	20.00	ng	0.00
38) Acenaphthene-d10	14.61	164	134309	20.00	ng	0.00
63) Phenanthrene-d10	17.35	188	442174	20.00	ng	0.00
75) Chrysene-d12	21.50	240	754521	20.00	ng	0.00
86) Perylene-d12	23.77	264	995126	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	109671	35.33	ng	0.00
7) Phenol-d6	7.15	99	209945	51.12	ng	0.00
23) Nitrobenzene-d5	9.16	82	189407	41.33	ng	0.00
41) 2,4,6-Tribromophenol	16.11	330	82585	22.00	ng	0.00
44) 2-Fluorobiphenyl	13.22	172	453113	32.43	ng	0.00
78) Terphenyl-d14	19.94	244	1285795	36.07	ng	0.00
Target Compounds						
49) Dimethylphthalate	14.07	163	122955	9.529	ng	Qvalue # 97
74) Fluoranthene	19.41	202	154172	4.601	ng	93
77) Pyrene	19.77	202	147355	3.594	ng	98
80) Benzo(a)anthracene	21.49	228	107351	2.420	ng	98
82) Chrysene	21.54	228	93234	2.203	ng	96
87) Benzo(b)fluoranthene	23.10	252	173623	3.027	ng	# 90
89) Benzo(a)pyrene	23.69	252	113579	2.035	ng	# 83

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

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