

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121918\
 Data File : BM018309.D
 Acq On : 19 Dec 2018 22:48
 Operator : JU/SJ
 Sample : J6378-22
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS002-0006-01

Quant Time: Dec 20 02:47:06 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 15 21:39:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	118079	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	427237	20.00	ng	-0.02
39) Acenaphthene-d10	14.14	164	205012	20.00	ng	-0.02
64) Phenanthrene-d10	16.90	188	412103	20.00	ng	-0.02
76) Chrysene-d12	21.13	240	491795	20.00	ng	-0.02
87) Perylene-d12	23.29	264	584190	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	776600	109.87	ng	0.00
7) Phenol-d6	6.69	99	1000635	101.85	ng	-0.02
23) Nitrobenzene-d5	8.64	82	642925	77.19	ng	-0.02
42) 2,4,6-Tribromophenol	15.65	330	284551	94.57	ng	-0.02
45) 2-Fluorobiphenyl	12.75	172	954540	60.31	ng	-0.02
79) Terphenyl-d14	19.56	244	960666	44.18	ng	-0.02
Target Compounds						
10) Phenol	6.71	94	63734	6.125	ng	83
50) Dimethylphthalate	13.62	163	79952	4.691	ng	99
83) Chrysene	21.16	228	62902	2.276	ng	96
86) Indeno(1,2,3-cd)pyrene	25.40	276	99534	3.616	ng	# 94
88) Benzo(b)fluoranthene	22.65	252	161736	4.929	ng	95
90) Benzo(a)pyrene	23.19	252	95320	3.054	ng	95
92) Benzo(g,h,i)perylene	26.06	276	96397	3.527	ng	93

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

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