

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM010919\
 Data File : BM018480.D
 Acq On : 09 Jan 2019 19:30
 Operator : JU/SJ
 Sample : K1006-17 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JC-03-010819-I

Manual Integrations
 APPROVED

Sohil
 1/11/2019 2:22:19 PM

Quant Time: Jan 10 06:18:31 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 09 12:39:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	592363	20.00	ng	0.00
21) Naphthalene-d8	10.55	136	2500161	20.00	ng	0.00
39) Acenaphthene-d10	14.41	164	1462515	20.00	ng	0.00
64) Phenanthrene-d10	17.16	188	3458768	20.00	ng	0.00
76) Chrysene-d12	21.36	240	2928126	20.00	ng	0.00
87) Perylene-d12	23.65	264	2677589	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1498131	46.70	ng	0.00
7) Phenol-d6	6.93	99	1966842	48.27	ng	0.00
23) Nitrobenzene-d5	8.92	82	1198628	27.79	ng	0.00
42) 2,4,6-Tribromophenol	15.90	330	848314	45.55	ng	0.00
45) 2-Fluorobiphenyl	13.02	172	2831950	25.27	ng	0.00
79) Terphenyl-d14	19.80	244	4377865	31.52	ng	0.01
Target Compounds						
10) Phenol	6.96	94	88915	2.147	ng	94
50) Dimethylphthalate	13.87	163	314888	2.622	ng	100
71) Phenanthrene	17.20	178	1967581	10.695	ng	99
75) Fluoranthene	19.22	202	1881359	8.870	ng	98
78) Pyrene	19.59	202	2107853	12.145	ng	99
81) Benzo(a)anthracene	21.34	228	769168	4.459	ng	96
83) Chrysene	21.39	228	717654	4.308	ng	97
88) Benzo(b)fluoranthene	22.96	252	655782m	4.314	ng	
90) Benzo(a)pyrene	23.54	252	503843	3.462	ng	# 93
92) Benzo(g,h,i)perylene	26.67	276	292955	2.042	ng	96

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

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