

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050919\
 Data File : BM020223.D
 Acq On : 09 May 2019 18:39
 Operator : JU/SJ
 Sample : K2645-07 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PL-02-050719-D

Quant Time: May 10 05:25:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 03:58:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	104932	20.00	ng	-0.02
21) Naphthalene-d8	10.57	136	391126	20.00	ng	-0.02
39) Acenaphthene-d10	14.42	164	236130	20.00	ng	-0.01
64) Phenanthrene-d10	17.17	188	554267	20.00	ng	-0.01
76) Chrysene-d12	21.36	240	629745	20.00	ng	-0.01
87) Perylene-d12	23.64	264	789939	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	152381	23.74	ng	-0.01
7) Phenol-d6	6.94	99	204802	23.35	ng	-0.01
23) Nitrobenzene-d5	8.95	82	140744	14.21	ng	-0.01
42) 2,4,6-Tribromophenol	15.91	330	81715	22.29	ng	-0.01
45) 2-Fluorobiphenyl	13.04	172	271022	14.12	ng	-0.02
79) Terphenyl-d14	19.79	244	467426	12.95	ng	-0.02
Target Compounds				Qvalue		
49) Acenaphthylene	14.15	152	131272	5.286	ng	99
58) Fluorene	15.47	166	58424	2.961	ng	95
71) Phenanthrene	17.21	178	581443	19.004	ng	100
72) Anthracene	17.30	178	225903	7.385	ng	99
75) Fluoranthene	19.23	202	1256316	32.453	ng	99
78) Pyrene	19.59	202	1162651	27.499	ng	99
81) Benzo(a)anthracene	21.34	228	748914	16.958	ng	97
83) Chrysene	21.39	228	642596	15.003	ng	98
86) Indeno(1,2,3-cd)pyrene	25.97	276	489396	9.606	ng	# 100
88) Benzo(b)fluoranthene	22.96	252	983225m	18.849	ng	
89) Benzo(k)fluoranthene	22.99	252	304261m	6.394	ng	
90) Benzo(a)pyrene	23.54	252	716012	15.112	ng	# 95
91) Dibenzo(a,h)anthracene	25.97	278	134857	2.737	ng	# 82
92) Benzo(g,h,i)perylene	26.69	276	466152	9.964	ng	98

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

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