

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM010919\
 Data File : BM018501.D
 Acq On : 10 Jan 2019 09:21
 Operator : JU/SJ
 Sample : K1044-05DL 10X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CPSB-10(10-15)DL

Manual Integrations
 APPROVED

Sohil
 1/10/2019 3:12:16 PM

Quant Time: Jan 10 11:05:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 10 05:57:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	653730	20.00	ng	0.00
21) Naphthalene-d8	10.55	136	2727370	20.00	ng	0.00
39) Acenaphthene-d10	14.41	164	1524114	20.00	ng	0.00
64) Phenanthrene-d10	17.16	188	3086809	20.00	ng	0.00
76) Chrysene-d12	21.36	240	2435432	20.00	ng	0.00
87) Perylene-d12	23.64	264	2487810	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	277804	7.85	ng	0.00
7) Phenol-d6	6.93	99	376787	8.38	ng	0.00
23) Nitrobenzene-d5	8.93	82	237960	5.06	ng	0.00
42) 2,4,6-Tribromophenol	15.90	330	134608	6.94	ng	0.00
45) 2-Fluorobiphenyl	13.02	172	616242	5.28	ng	0.00
79) Terphenyl-d14	19.79	244	688055	5.96	ng	0.00
Target Compounds						
52) Acenaphthene	14.47	154	305788	3.323	ng	98
58) Fluorene	15.46	166	506536	4.472	ng	97
71) Phenanthrene	17.20	178	3629493	22.106	ng	99
72) Anthracene	17.29	178	1081926	6.853	ng	99
75) Fluoranthene	19.22	202	3677282	19.426	ng	96
78) Pyrene	19.59	202	4810600	33.324	ng	99
81) Benzo(a)anthracene	21.34	228	1973639	13.756	ng	98
83) Chrysene	21.39	228	1859839	13.424	ng	98
86) Indeno(1,2,3-cd)pyrene	25.96	276	735469	4.604	ng	# 93
88) Benzo(b)fluoranthene	22.96	252	1676334m	11.869	ng	
89) Benzo(k)fluoranthene	22.99	252	509395m	3.579	ng	
90) Benzo(a)pyrene	23.54	252	1377386	10.187	ng	99
92) Benzo(a,h,i)perylene	26.67	276	734045	5.506	ng	97

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

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