

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022619\
 Data File : BM018942.D
 Acq On : 26 Feb 2019 22:50
 Operator : JU/SJ
 Sample : K1348-07 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OK-02-022519

Manual Integrations
 APPROVED

Sohil
 2/27/2019 4:20:25 PM

Quant Time: Feb 27 00:59:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 11 16:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	372794	20.00	ng	-0.02
21) Naphthalene-d8	10.48	136	1405524	20.00	ng	-0.02
39) Acenaphthene-d10	14.34	164	670054	20.00	ng	-0.02
64) Phenanthrene-d10	17.10	188	1314795	20.00	ng	-0.02
76) Chrysene-d12	21.29	240	1358218	20.00	ng	-0.02
87) Perylene-d12	23.54	264	1500252	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	1217161	53.50	ng	-0.02
7) Phenol-d6	6.87	99	1622769	50.63	ng	-0.02
23) Nitrobenzene-d5	8.86	82	1030856	33.91	ng	-0.03
42) 2,4,6-Tribromophenol	15.84	330	409377	42.38	ng	-0.02
45) 2-Fluorobiphenyl	12.96	172	1526613	30.24	ng	-0.02
79) Terphenyl-d14	19.73	244	1691100	25.68	ng	-0.02
Target Compounds						
71) Phenanthrene	17.14	178	333329	4.717	ng	Qvalue 100
75) Fluoranthene	19.16	202	661787	8.108	ng	99
78) Pyrene	19.53	202	598366	7.585	ng	99
81) Benzo(a)anthracene	21.27	228	333733	4.215	ng	96
83) Chrysene	21.33	228	301890	3.978	ng	96
84) Bis(2-ethylhexyl)phthalate	21.20	149	112589	2.140	ng	98
86) Indeno(1,2,3-cd)pyrene	25.82	276	235286	2.685	ng	96
88) Benzo(b)fluoranthene	22.86	252	443923m	5.172	ng	
89) Benzo(k)fluoranthene	22.90	252	172952m	2.024	ng	
90) Benzo(a)pyrene	23.44	252	332972	4.095	ng	# 93
92) Benzo(g,h,i)perylene	26.53	276	241708	3.223	ng	96

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

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