

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121918\
 Data File : BM018295.D
 Acq On : 19 Dec 2018 13:18
 Operator : JU/SJ
 Sample : J6319-11 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 065_SW-W-1

Manual Integrations
 APPROVED

Sohil
 12/20/2018 2:55:04 PM

Quant Time: Dec 19 14:33:48 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	7.49	152	93793	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	349567	20.00	ng	-0.02
39) Acenaphthene-d10	14.15	164	185256	20.00	ng	-0.02
64) Phenanthrene-d10	16.91	188	409568	20.00	ng	-0.01
76) Chrysene-d12	21.13	240	501265	20.00	ng	-0.02
87) Perylene-d12	23.29	264	560690	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	46602	8.30	ng	0.00
7) Phenol-d6	6.69	99	53575	6.87	ng	-0.01
23) Nitrobenzene-d5	8.66	82	31003	4.55	ng	0.00
42) 2,4,6-Tribromophenol	15.66	330	16228	5.97	ng	0.00
45) 2-Fluorobiphenyl	12.76	172	55778	3.90	ng	-0.01
79) Terphenyl-d14	19.56	244	87541	3.95	ng	-0.01
Target Compounds						
75) Fluoranthene	18.98	202	146305	5.656	ng	98
78) Pyrene	19.35	202	140891	4.717	ng	99
81) Benzo(a)anthracene	21.12	228	188867	6.450	ng	99
83) Chrysene	21.16	228	175224m	6.221	ng	
86) Indeno(1,2,3-cd)pyrene	25.41	276	300884	10.725	ng	96
88) Benzo(b)fluoranthene	22.65	252	413351m	13.124	ng	
89) Benzo(k)fluoranthene	22.69	252	121031m	3.859	ng	
90) Benzo(a)pyrene	23.19	252	372861	12.447	ng	99
91) Dibenzo(a,h)anthracene	25.41	278	100645	3.627	ng	# 78
92) Benzo(g,h,i)perylene	26.06	276	293157	11.177	ng	98

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

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