

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121918\
 Data File : BM018316.D
 Acq On : 20 Dec 2018 03:01
 Operator : JU/SJ
 Sample : J6379-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS003-0612-01

Manual Integrations
 APPROVED

Sohil
 12/20/2018 2:55:18 PM

Quant Time: Dec 20 07:23:55 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	7.48	152	95420	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	345192	20.00	ng	-0.02
39) Acenaphthene-d10	14.15	164	176020	20.00	ng	-0.02
64) Phenanthrene-d10	16.90	188	362217	20.00	ng	-0.02
76) Chrysene-d12	21.13	240	436476	20.00	ng	-0.02
87) Perylene-d12	23.29	264	489169	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	635414	111.24	ng	0.00
7) Phenol-d6	6.69	99	829821	104.52	ng	-0.02
23) Nitrobenzene-d5	8.64	82	540002	80.24	ng	-0.02
42) 2,4,6-Tribromophenol	15.65	330	257589	99.71	ng	-0.02
45) 2-Fluorobiphenyl	12.75	172	1013173	74.55	ng	-0.02
79) Terphenyl-d14	19.56	244	1275656	66.10	ng	-0.01
Target Compounds						
10) Phenol	6.71	94	48290	5.743	ng	88
50) Dimethylphthalate	13.62	163	56335	3.850	ng	# 98
75) Fluoranthene	18.98	202	207484	9.070	ng	99
78) Pyrene	19.34	202	178151	6.850	ng	99
81) Benzo(a)anthracene	21.11	228	190597	7.475	ng	98
83) Chrysene	21.16	228	211072	8.607	ng	99
86) Indeno(1,2,3-cd)pyrene	25.41	276	155535	6.367	ng	# 93
88) Benzo(b)fluoranthene	22.64	252	326749	11.891	ng	100
89) Benzo(k)fluoranthene	22.69	252	105513m	3.856	ng	
90) Benzo(a)pyrene	23.19	252	187602	7.179	ng	98
92) Benzo(g,h,i)perylene	26.06	276	138575	6.056	ng	98

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

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