

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120518\
 Data File : BM017932.D
 Acq On : 05 Dec 2018 09:36
 Operator : JU/SJ
 Sample : J6162-02
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A0C10-01B1-(0-6)

Quant Time: Dec 05 11:11:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM111218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 12 18:37:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	141462	20.00	ng	-0.04
21) Naphthalene-d8	10.35	136	556521	20.00	ng	-0.04
39) Acenaphthene-d10	14.23	164	308461	20.00	ng	-0.04
64) Phenanthrene-d10	16.99	188	694748	20.00	ng	-0.03
76) Chrysene-d12	21.19	240	761908	20.00	ng	-0.02
87) Perylene-d12	23.37	264	800819	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.21	112	926197	110.50	ng	-0.03
7) Phenol-d6	6.77	99	1194861	102.06	ng	-0.03
23) Nitrobenzene-d5	8.73	82	735076	68.19	ng	-0.04
42) 2,4,6-Tribromophenol	15.73	330	505418	114.03	ng	-0.03
45) 2-Fluorobiphenyl	12.84	172	1745048	73.39	ng	-0.04
79) Terphenyl-d14	19.63	244	2590070	77.05	ng	-0.02
Target Compounds						
10) Phenol	6.80	94	60786	4.947	ng	Qvalue 94
50) Dimethylphthalate	13.70	163	115675	4.612	ng	99
75) Fluoranthene	19.06	202	231944	5.438	ng	99
78) Pyrene	19.42	202	283975	6.041	ng	100
81) Benzo(a)anthracene	21.18	228	137273	3.127	ng	94
83) Chrysene	21.23	228	141099	3.310	ng	97
84) Bis(2-ethylhexyl)phthalate	21.12	149	833310	29.522	ng	100
86) Indeno(1,2,3-cd)pyrene	25.55	276	130942	3.843	ng	97
88) Benzo(b)fluoranthene	22.73	252	210222	4.482	ng	# 93
90) Benzo(a)pyrene	23.28	252	170570	3.989	ng	96
92) Benzo(g,h,i)perylene	26.21	276	144915	4.287	ng	96

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

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