

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF121917\
 Data File : BF101524.D
 Acq On : 19 Dec 2017 16:05
 Operator : SJ/JU
 Sample : I6908-09
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 5

Manual Integrations
 APPROVED

Sohil
 12/20/2017 3:47:48 PM

Quant Time: Dec 19 16:59:59 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 18 15:38:59 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	126074	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	508614	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	220479	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	375368	20.00	ng	0.00
75) Chrysene-d12	13.99	240	225771	20.00	ng	0.00
86) Perylene-d12	15.46	264	270191	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	855532	101.08	ng	0.01
7) Phenol-d6	6.45	99	1029129	97.70	ng	0.00
23) Nitrobenzene-d5	7.37	82	653527	71.53	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	209461	98.59	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	1076848	75.76	ng	0.00
78) Terphenyl-d14	12.92	244	800946	85.52	ng	0.00
Target Compounds						
10) Phenol	6.46	94	47423	3.950	ng	81
49) Dimethylphthalate	9.56	163	52837	2.979	ng	100
70) Phenanthrene	11.36	178	106686	5.111	ng	97
74) Fluoranthene	12.56	202	167276	7.634	ng	96
77) Pyrene	12.79	202	120611	7.118	ng	98
80) Benzo(a)anthracene	13.97	228	61660	4.136	ng	98
82) Chrysene	14.02	228	68144	4.841	ng	97
85) Indeno(1,2,3-cd)pyrene	16.94	276	39272	3.133	ng	96
87) Benzo(b)fluoranthene	15.03	252	83173m	5.050	ng	
89) Benzo(a)pyrene	15.40	252	52161	3.436	ng	98
91) Benzo(g,h,i)perylene	17.40	276	36207m	2.805	ng	

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

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