

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF120117\
 Data File : BF101062.D
 Acq On : 1 Dec 2017 14:50
 Operator : SJ/JU
 Sample : I6610-03 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-3(10-12)-20171127

Manual Integrations
 APPROVED

Sohil
 12/4/2017 11:56:32 AM

Quant Time: Dec 04 00:15:33 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 01 15:52:19 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	51489	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	202257	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	91378	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	174482	20.00	ng	0.00
75) Chrysene-d12	14.02	240	112409	20.00	ng	0.00
86) Perylene-d12	15.49	264	121577	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	66708	21.41	ng	-0.01
7) Phenol-d6	6.47	99	81242	21.48	ng	-0.01
23) Nitrobenzene-d5	7.40	82	47224	13.93	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	13592	12.38	ng	-0.01
44) 2-Fluorobiphenyl	9.20	172	108504	16.25	ng	-0.01
78) Terphenyl-d14	12.95	244	88971	17.31	ng	-0.01
Target Compounds						
70) Phenanthrene	11.39	178	46118	4.800	ng	98
74) Fluoranthene	12.59	202	76634	7.195	ng	94
77) Pyrene	12.82	202	66691	8.598	ng	98
79) Butylbenzylphthalate	13.43	149	23483	6.867	ng	87
80) Benzo(a)anthracene	14.00	228	29452	4.150	ng	93
82) Chrysene	14.04	228	30000m	4.551	ng	
85) Indeno(1,2,3-cd)pyrene	16.96	276	18289	3.121	ng	# 91
87) Benzo(b)fluoranthene	15.05	252	40827m	5.606	ng	
89) Benzo(a)pyrene	15.42	252	28605	4.321	ng	95
91) Benzo(g,h,i)perylene	17.41	276	22143	4.026	ng	94

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

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