

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF120117\
 Data File : BF101066.D
 Acq On : 1 Dec 2017 16:37
 Operator : SJ/JU
 Sample : I6618-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ENV-106-6(0-5)

Manual Integrations
 APPROVED

Sohil
 12/4/2017 11:56:43 AM

Quant Time: Dec 04 00:35:58 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	54900	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	213493	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	95996	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	164725	20.00	ng	0.00
75) Chrysene-d12	14.02	240	128045	20.00	ng	0.00
86) Perylene-d12	15.50	264	119115	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	313388	94.33	ng	0.00
7) Phenol-d6	6.48	99	379955	94.20	ng	0.00
23) Nitrobenzene-d5	7.41	82	234754	65.59	ng	0.00
41) 2,4,6-Tribromophenol	10.67	330	72591	62.95	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	460448	65.63	ng	0.00
78) Terphenyl-d14	12.96	244	350284	59.84	ng	0.00
Target Compounds						
10) Phenol	6.49	94	15716	3.504	ng	97
49) Dimethylphthalate	9.59	163	23228	3.000	ng	# 96
70) Phenanthrene	11.40	178	76023	8.381	ng	97
71) Anthracene	11.44	178	23488	2.558	ng	97
74) Fluoranthene	12.59	202	145204	14.440	ng	96
77) Pyrene	12.82	202	134079	15.175	ng	98
80) Benzo(a)anthracene	14.01	228	71383	8.830	ng	98
82) Chrysene	14.04	228	68778	9.160	ng	99
85) Indeno(1,2,3-cd)pyrene	16.97	276	30353	4.547	ng	95
87) Benzo(b)fluoranthene	15.06	252	85850m	12.033	ng	
88) Benzo(k)fluoranthene	15.09	252	23477m	3.340	ng	
89) Benzo(a)pyrene	15.43	252	57844	8.918	ng	95
91) Benzo(a,h,i)perylene	17.43	276	27180m	5.044	ng	

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

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