

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110117\
 Data File : BF100050.D
 Acq On : 1 Nov 2017 18:14
 Operator : SJ/JU
 Sample : I6200-01 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-01-102717

Manual Integrations
 APPROVED

Sohil
 11/2/2017 4:27:10 PM

Quant Time: Nov 01 18:56:40 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 18:13:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	83966	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	340243	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	148100	20.00	ng	0.00
63) Phenanthrene-d10	11.31	188	234595	20.00	ng	0.00
75) Chrysene-d12	13.96	240	160984	20.00	ng	-0.01
86) Perylene-d12	15.44	264	163201	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	101845	19.04	ng	0.01
7) Phenol-d6	6.43	99	129640	20.44	ng	0.00
23) Nitrobenzene-d5	7.35	82	68994	13.05	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	24631	18.25	ng	0.00
44) 2-Fluorobiphenyl	9.13	172	152360	15.87	ng	0.00
78) Terphenyl-d14	12.89	244	105150	14.27	ng	-0.01
Target Compounds						
70) Phenanthrene	11.33	178	72290	5.460	ng	98
74) Fluoranthene	12.53	202	94798	6.890	ng	99
77) Pyrene	12.76	202	86322	7.052	ng	98
80) Benzo(a)anthracene	13.95	228	38665	3.789	ng	98
82) Chrysene	13.99	228	35227	3.672	ng	96
85) Indeno(1,2,3-cd)pyrene	16.94	276	19937	2.264	ng	# 86
87) Benzo(b)fluoranthene	15.02	252	44141m	4.283	ng	
89) Benzo(a)pyrene	15.39	252	31583	3.412	ng	# 90
91) Benzo(g,h,i)perylene	17.39	276	19439	2.416	ng	# 91

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

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