

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF022818\
 Data File : BF103301.D
 Acq On : 28 Feb 2018 18:57
 Operator : SJ/JU
 Sample : J1710-04 10X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-201

Manual Integrations
 APPROVED

Sohil
 3/1/2018 11:43:11 AM

Quant Time: Mar 01 00:46:03 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 17:05:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	153365	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	647357	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	276456	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	434818	20.00	ng	0.00
75) Chrysene-d12	14.06	240	281972	20.00	ng	0.00
86) Perylene-d12	15.56	264	262088	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	129495	12.77	ng	0.00
7) Phenol-d6	6.50	99	164355	13.25	ng	0.00
23) Nitrobenzene-d5	7.43	82	97275	8.58	ng	-0.01
41) 2,4,6-Tribromophenol	10.71	330	23143	9.89	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	163541	6.10	ng	0.00
78) Terphenyl-d14	12.99	244	106001	9.04	ng	0.00
Target Compounds						
70) Phenanthrene	11.43	178	179202	7.489	ng	98
74) Fluoranthene	12.62	202	168918	7.025	ng	98
77) Pyrene	12.85	202	174323	7.929	ng	99
80) Benzo(a)anthracene	14.04	228	60288	3.295	ng	98
82) Chrysene	14.08	228	61553	3.465	ng	98
87) Benzo(b)fluoranthene	15.12	252	64569m	3.747	ng	
89) Benzo(a)pyrene	15.49	252	47777	3.105	ng	# 93
91) Benzo(g,h,i)perylene	17.54	276	23892	2.056	ng	# 90

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

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