

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110117\
 Data File : BF100049.D
 Acq On : 1 Nov 2017 17:48
 Operator : SJ/JU
 Sample : I6206-05 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-102717-C

Manual Integrations
 APPROVED

Sohil
 11/2/2017 4:27:07 PM

Quant Time: Nov 01 18:46:38 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	86640	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	343597	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	149876	20.00	ng	0.00
63) Phenanthrene-d10	11.31	188	236490	20.00	ng	0.00
75) Chrysene-d12	13.97	240	155538	20.00	ng	0.00
86) Perylene-d12	15.46	264	155159	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	52329	9.48	ng	0.02
7) Phenol-d6	6.44	99	68579	10.48	ng	0.01
23) Nitrobenzene-d5	7.36	82	41493	7.77	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	13725	10.05	ng	0.00
44) 2-Fluorobiphenyl	9.13	172	87713	9.03	ng	0.00
78) Terphenyl-d14	12.90	244	58182	8.17	ng	0.00
Target Compounds						
74) Fluoranthene	12.54	202	35687	2.573	ng	93
77) Pyrene	12.77	202	39266	3.320	ng	97
80) Benzo(a)anthracene	13.96	228	23242	2.357	ng	95
82) Chrysene	14.00	228	22650	2.443	ng	98
87) Benzo(b)fluoranthene	15.03	252	29400m	3.001	ng	
89) Benzo(a)pyrene	15.40	252	21191	2.408	ng	# 91

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

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