

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF022818\
 Data File : BF103300.D
 Acq On : 28 Feb 2018 18:30
 Operator : SJ/JU
 Sample : J1710-07 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-215

Manual Integrations
 APPROVED

Sohil
 3/1/2018 11:43:09 AM

Quant Time: Mar 01 00:43:01 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	147170	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	595920	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	238085	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	337650	20.00	ng	0.00
75) Chrysene-d12	14.06	240	259707	20.00	ng	0.00
86) Perylene-d12	15.56	264	253179	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	99299	10.20	ng	0.00
7) Phenol-d6	6.50	99	126959	10.66	ng	0.00
23) Nitrobenzene-d5	7.44	82	78920	7.56	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	14923	7.41	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	120019	4.78	ng	0.00
78) Terphenyl-d14	12.99	244	69791	6.46	ng	0.00
Target Compounds						
70) Phenanthrene	11.43	178	118407	6.372	ng	98
74) Fluoranthene	12.62	202	124632	6.675	ng	99
77) Pyrene	12.85	202	110075	5.436	ng	99
80) Benzo(a)anthracene	14.04	228	50781	3.014	ng	96
82) Chrysene	14.08	228	49763	3.041	ng	97
87) Benzo(b)fluoranthene	15.12	252	58015m	3.485	ng	
89) Benzo(a)pyrene	15.49	252	40295	2.711	ng	# 84

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

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