

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111618\
 Data File : BF110828.D
 Acq On : 16 Nov 2018 20:28
 Operator : JU/SJ
 Sample : J5951-01 10X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FDSB-02(10-12)

Manual Integrations
 APPROVED

Sohil
 11/19/2018 11:56:03 AM

Quant Time: Nov 17 00:09:21 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	54626	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	225842	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	106475	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	176720	20.00	ng	-0.01
76) Chrysene-d12	14.13	240	151012	20.00	ng	-0.01
87) Perylene-d12	15.66	264	113224	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	28745	8.55	ng	0.00
7) Phenol-d6	6.57	99	37551	8.73	ng	0.00
23) Nitrobenzene-d5	7.51	82	21052m	5.37	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	5988	5.58	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	48095	6.45	ng	0.00
79) Terphenyl-d14	13.06	244	38420	4.76	ng	-0.01
Target Compounds						
37) 2-Methylnaphthalene	8.95	142	713947	102.724	ng	Qvalue 100
38) 1-Methylnaphthalene	9.05	142	425671m	63.685	ng	
55) Dibenzofuran	10.19	168	27879	2.926	ng	# 65
58) Fluorene	10.54	166	29966	4.094	ng	# 66
71) Phenanthrene	11.50	178	25840	2.729	ng	97

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

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