

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
 Data File : BF109701.D
 Acq On : 9 Oct 2018 18:07
 Operator : JU/SJ
 Sample : J5325-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SOIL-DEBRI

Manual Integrations
 APPROVED

Sohil
 10/10/2018 11:37:05 AM

Quant Time: Oct 10 09:30:32 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF100818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 08 16:02:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	93993	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	375928	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	174820	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	300263	20.00	ng	0.00
75) Chrysene-d12	13.83	240	204269	20.00	ng	0.00
86) Perylene-d12	15.23	264	204306	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	571098	107.85	ng	0.00
7) Phenol-d6	6.35	99	756144	106.89	ng	-0.01
23) Nitrobenzene-d5	7.26	82	518899	76.09	ng	-0.01
41) 2,4,6-Tribromophenol	10.51	330	191717	100.65	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	971311	76.52	ng	0.00
78) Terphenyl-d14	12.78	244	698242	66.17	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.45	163	119072	9.287	ng	99
74) Fluoranthene	12.41	202	50769	3.234	ng	98
77) Pyrene	12.63	202	43596	2.913	ng	98
82) Chrysene	13.86	228	27435	2.317	ng	96
85) Indeno(1,2,3-cd)pyrene	16.59	276	18322	2.162	ng	# 91
87) Benzo(b)fluoranthene	14.84	252	34577m	3.063	ng	
89) Benzo(a)pyrene	15.17	252	23584	2.302	ng	# 70
91) Benzo(g,h,i)perylene	16.99	276	17719	2.109	ng	# 88

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

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