

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092218\
 Data File : BF109253.D
 Acq On : 23 Sep 2018 6:27
 Operator : JU/SJ
 Sample : J5017-01
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RA-091918-N-NW-055

Manual Integrations
 APPROVED

Sohil
 9/24/2018 3:53:38 PM

Quant Time: Sep 24 11:31:03 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	115741	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	431825	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	185631	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	263680	20.00	ng	0.00
75) Chrysene-d12	13.85	240	187425	20.00	ng	0.00
86) Perylene-d12	15.26	264	162899	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	638695	90.89	ng	0.00
7) Phenol-d6	6.35	99	811455	90.50	ng	-0.01
23) Nitrobenzene-d5	7.27	82	507395	69.36	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	154614	84.49	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	795084	64.60	ng	-0.01
78) Terphenyl-d14	12.81	244	416081	54.48	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.46	163	255227	18.903	ng	# 99
70) Phenanthrene	11.24	178	28124	2.136	ng	# 80
77) Pyrene	12.66	202	30190	2.248	ng	# 79
82) Chrysene	13.87	228	35452	3.168	ng	# 50
87) Benzo(b)fluoranthene	14.86	252	21580m	2.411	ng	

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

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