

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102618\
 Data File : BF110214.D
 Acq On : 26 Oct 2018 17:17
 Operator : JU/SJ
 Sample : J5658-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-TS-001

Manual Integrations
 APPROVED

Sohil
 10/29/2018 11:56:40 AM

Quant Time: Oct 27 00:07:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 23 15:06:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	131710	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	509920	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	209225	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	319024	20.00	ng	-0.01
76) Chrysene-d12	14.15	240	237103	20.00	ng	-0.01
87) Perylene-d12	15.68	264	164162	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	748879	92.59	ng	0.00
7) Phenol-d6	6.59	99	935764	90.45	ng	-0.01
23) Nitrobenzene-d5	7.53	82	571002	66.42	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	163168	78.73	ng	-0.01
45) 2-Fluorobiphenyl	9.33	172	901587	67.67	ng	-0.01
79) Terphenyl-d14	13.09	244	610301	53.53	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	106913	9.668	ng	Qvalue # 60
50) Dimethylphthalate	9.72	163	142572	9.231	ng	99
75) Fluoranthene	12.72	202	44971	2.655	ng	98
78) Pyrene	12.95	202	39515	2.325	ng	96
88) Benzo(b)fluoranthene	15.22	252	29495m	3.111	ng	
90) Benzo(a)pyrene	15.61	252	18830	2.159	ng	# 92

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

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