

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
 Data File : BF114141.D
 Acq On : 11 May 2019 00:33
 Operator : JU/SJ
 Sample : K2761-03 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-SF-03-006-B

Manual Integrations
 APPROVED

Sohil
 5/13/2019 8:34:31 PM

Quant Time: May 11 04:54:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 10 15:56:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	334096	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	1222029	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	512662	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	965761	20.00	ng	0.00
76) Chrysene-d12	14.01	240	626478	20.00	ng	-0.01
87) Perylene-d12	15.48	264	514244	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	251472	12.11	ng	0.00
7) Phenol-d6	6.49	99	542066	22.08	ng	0.00
23) Nitrobenzene-d5	7.41	82	296702	13.63	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	30139	5.69	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	566853	16.73	ng	0.00
79) Terphenyl-d14	12.95	244	550049	18.10	ng	0.00
Target Compounds						
71) Phenanthrene	11.39	178	192698	4.101	ng	Qvalue 96
75) Fluoranthene	12.58	202	210297	4.156	ng	95
78) Pyrene	12.81	202	205735	4.082	ng	97
81) Benzo(a)anthracene	14.00	228	81815	2.019	ng	# 89
88) Benzo(b)fluoranthene	15.05	252	68480m	2.079	ng	

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

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