

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
 Data File : BF114140.D
 Acq On : 11 May 2019 00:06
 Operator : JU/SJ
 Sample : K2761-01 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 5/13/2019 8:34:29 PM

Quant Time: May 11 04:48:30 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	6.85	152	322459	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	1197592	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	495038	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	940329	20.00	ng	0.00
76) Chrysene-d12	14.01	240	617832	20.00	ng	-0.01
87) Perylene-d12	15.47	264	505356	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	169246	8.45	ng	0.00
7) Phenol-d6	6.48	99	570415	24.07	ng	-0.01
23) Nitrobenzene-d5	7.41	82	326612	15.31	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	12274	2.40	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	588395	17.98	ng	0.00
79) Terphenyl-d14	12.95	244	556995	18.58	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.60	163	86230	2.373	ng	99
71) Phenanthrene	11.39	178	321051	7.018	ng	98
75) Fluoranthene	12.58	202	441178	8.955	ng	98
78) Pyrene	12.81	202	413849	8.327	ng	98
81) Benzo(a)anthracene	14.00	228	183053	4.581	ng	93
83) Chrysene	14.03	228	163237	4.167	ng	97
88) Benzo(b)fluoranthene	15.05	252	147842m	4.566	ng	
90) Benzo(a)pyrene	15.41	252	98703	3.464	ng	# 81
92) Benzo(g,h,i)perylene	17.39	276	57831	2.437	ng	# 83

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

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