

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050119\
 Data File : BF114067.D
 Acq On : 1 May 2019 14:21
 Operator : JU/SJ
 Sample : K2194-17RE 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-03-042919-IRE

Manual Integrations
 APPROVED

Sohil
 5/2/2019 4:53:46 PM

Quant Time: May 02 04:15:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 23 03:32:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	93712	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	357863	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	177607	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	349598	20.00	ng	0.00
76) Chrysene-d12	14.08	240	267754	20.00	ng	0.01
87) Perylene-d12	15.60	264	194451	20.00	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	227676	41.55	ng	0.00
7) Phenol-d6	6.52	99	306958	44.65	ng	0.00
23) Nitrobenzene-d5	7.45	82	190714	32.90	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	97563	45.09	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	369493	32.54	ng	0.00
79) Terphenyl-d14	13.00	244	400054	30.63	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.65	163	54548	4.247	ng	99
71) Phenanthrene	11.44	178	62014	3.530	ng	98
75) Fluoranthene	12.64	202	98187	5.122	ng	97
78) Pyrene	12.87	202	93938	5.017	ng	98
81) Benzo(a)anthracene	14.07	228	49201	3.119	ng	# 91
83) Chrysene	14.10	228	46127	3.002	ng	97
88) Benzo(b)fluoranthene	15.16	252	50412m	4.098	ng	
90) Benzo(a)pyrene	15.53	252	29323	2.756	ng	# 72

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

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