

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050119\
 Data File : BF114066.D
 Acq On : 1 May 2019 13:55
 Operator : JU/SJ
 Sample : K2194-13RE 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 5/2/2019 4:53:43 PM

Quant Time: May 02 03:35:04 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	96266	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	324026	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	202263	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	320007	20.00	ng	0.00
76) Chrysene-d12	14.07	240	246594	20.00	ng	0.00
87) Perylene-d12	15.57	264	151946	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	270850	48.12	ng	0.00
7) Phenol-d6	6.52	99	436438	61.81	ng	0.00
23) Nitrobenzene-d5	7.45	82	207683	39.57	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	101655	41.26	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	421375	32.58	ng	0.00
79) Terphenyl-d14	13.01	244	457684	38.05	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	9.64	163	60026	4.103	ng	99
71) Phenanthrene	11.44	178	52121	3.241	ng	98
75) Fluoranthene	12.64	202	75683	4.313	ng	96
78) Pyrene	12.87	202	78707	4.564	ng	97
81) Benzo(a)anthracene	14.06	228	36128	2.487	ng	# 85
83) Chrysene	14.10	228	37782	2.670	ng	# 90
88) Benzo(b)fluoranthene	15.14	252	33867m	3.523	ng	
90) Benzo(a)pyrene	15.51	252	18818	2.263	ng	# 42

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

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