

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040119\
 Data File : BF113474.D
 Acq On : 1 Apr 2019 18:50
 Operator : JU/SJ
 Sample : K2184-01 4X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-0D015-03-001

Manual Integrations
 APPROVED

Sohil
 4/2/2019 3:49:32 PM

Quant Time: Apr 02 04:16:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	142814	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	555130	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	248228	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	427488	20.00	ng	0.00
76) Chrysene-d12	13.84	240	380026	20.00	ng	0.00
87) Perylene-d12	15.26	264	265361	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	89131	10.20	ng	0.00
7) Phenol-d6	6.35	99	190717	17.26	ng	0.00
23) Nitrobenzene-d5	7.27	82	105685	11.30	ng	0.00
42) 2,4,6-Tribromophenol	10.53	330	14138	4.61	ng	0.00
45) 2-Fluorobiphenyl	0.00	172	0d	0.00	ng	0.00
79) Terphenyl-d14	12.79	244	230648	13.38	ng	0.00
Target Compounds						
71) Phenanthrene	11.24	178	121341	5.716	ng	97
75) Fluoranthene	12.42	202	142609	5.942	ng	99
78) Pyrene	12.65	202	133981	5.088	ng	99
81) Benzo(a)anthracene	13.83	228	63572	2.922	ng	# 91
83) Chrysene	13.87	228	65720	2.974	ng	# 95
88) Benzo(b)fluoranthene	14.86	252	54360m	3.298	ng	
90) Benzo(a)pyrene	15.20	252	36616	2.506	ng	# 73

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

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