

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032819\
 Data File : BF113405.D
 Acq On : 29 Mar 2019 7:27
 Operator : JU/SJ
 Sample : K1950-03RE 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-SF-03-005-BRE

Manual Integrations
 APPROVED

Sohil
 3/29/2019 4:47:29 PM

Quant Time: Mar 29 08:43:46 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	178282	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	689840	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	308390	20.00	ng	0.00
64) Phenanthrene-d10	11.21	188	544441	20.00	ng	0.00
76) Chrysene-d12	13.84	240	452015	20.00	ng	0.00
87) Perylene-d12	15.24	264	359211	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	201915	18.52	ng	0.00
7) Phenol-d6	6.35	99	363781	26.37	ng	-0.01
23) Nitrobenzene-d5	7.26	82	201797	17.36	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	28148	7.39	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	470744	7.93	ng	0.00
79) Terphenyl-d14	12.79	244	415818	20.28	ng	-0.01
Target Compounds						
71) Phenanthrene	11.23	178	212612	7.864	ng	98
75) Fluoranthene	12.42	202	214335	7.012	ng	98
78) Pyrene	12.64	202	216019	6.896	ng	98
81) Benzo(a)anthracene	13.83	228	96580	3.733	ng	# 92
83) Chrysene	13.86	228	99459	3.784	ng	# 95
88) Benzo(b)fluoranthene	14.85	252	91360m	4.094	ng	
90) Benzo(a)pyrene	15.19	252	60029	3.034	ng	# 78

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

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