

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032919\
 Data File : BF113430.D
 Acq On : 29 Mar 2019 23:15
 Operator : JU/SJ
 Sample : K1698-03 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-05-032819

Manual Integrations
 APPROVED

Sohil
 4/1/2019 5:46:38 PM

Quant Time: Apr 01 08:31:34 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	117601	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	469728	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	231673	20.00	ng	0.00
64) Phenanthrene-d10	11.21	188	451062	20.00	ng	0.00
76) Chrysene-d12	13.84	240	442492	20.00	ng	0.00
87) Perylene-d12	15.24	264	350517	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	133616	18.58	ng	0.00
7) Phenol-d6	6.35	99	192370	21.14	ng	-0.01
23) Nitrobenzene-d5	7.26	82	112090	14.16	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	55233	19.30	ng	0.00
45) 2-Fluorobiphenyl	0.00	172	0d	0.00	ng	
79) Terphenyl-d14	12.79	244	283973	14.15	ng	0.00
Target Compounds						
71) Phenanthrene	11.23	178	101891	4.549	ng	98
75) Fluoranthene	12.42	202	219023	8.649	ng	99
78) Pyrene	12.64	202	219101	7.145	ng	98
81) Benzo(a)anthracene	13.83	228	99466	3.927	ng	99
83) Chrysene	13.86	228	92455	3.594	ng	98
88) Benzo(b)fluoranthene	14.85	252	108320m	4.975	ng	
90) Benzo(a)pyrene	15.19	252	68573	3.552	ng	# 90

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

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