

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032619\
 Data File : BF113315.D
 Acq On : 27 Mar 2019 2:46
 Operator : JU/SJ
 Sample : K2078-01RE 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-SF-01-005-ARE

Manual Integrations
 APPROVED

Sohil
 3/27/2019 4:49:56 PM

Quant Time: Mar 27 08:28:51 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	160582	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	590904	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	250345	20.00	ng	0.00
64) Phenanthrene-d10	11.21	188	410892	20.00	ng	0.00
76) Chrysene-d12	13.83	240	405724	20.00	ng	-0.01
87) Perylene-d12	15.24	264	402565	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	83044	8.46	ng	0.00
7) Phenol-d6	6.34	99	421676	33.94	ng	-0.02
23) Nitrobenzene-d5	7.26	82	330339	33.18	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	12969	4.19	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	620959	26.80	ng	0.00
79) Terphenyl-d14	12.79	244	523147	28.42	ng	-0.01
Target Compounds						
71) Phenanthrene	11.23	178	241683	11.844	ng	Qvalue 98
72) Anthracene	11.29	178	60579	2.923	ng	98
75) Fluoranthene	12.42	202	280938	12.179	ng	99
78) Pyrene	12.64	202	252924	8.996	ng	98
81) Benzo(a)anthracene	13.83	228	131315	5.654	ng	98
83) Chrysene	13.86	228	126518	5.363	ng	97
86) Indeno(1,2,3-cd)pyrene	16.59	276	55176	2.725	ng	98
88) Benzo(b)fluoranthene	14.84	252	158676m	6.345	ng	
90) Benzo(a)pyrene	15.18	252	103582	4.672	ng	# 93
92) Benzo(g,h,i)perylene	17.00	276	52001	2.690	ng	# 94

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

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