

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032119\
 Data File : BF113218.D
 Acq On : 21 Mar 2019 19:48
 Operator : JU/SJ
 Sample : K2004-12 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-23(4-5)

Manual Integrations
 APPROVED

Sohil
 3/22/2019 3:36:02 PM

Quant Time: Mar 22 01:52:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 13 03:42:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	166683	20.00	ng	-0.01
21) Naphthalene-d8	8.00	136	677023	20.00	ng	-0.01
39) Acenaphthene-d10	9.76	164	260952	20.00	ng	-0.01
64) Phenanthrene-d10	11.23	188	517564	20.00	ng	-0.01
76) Chrysene-d12	13.87	240	357513	20.00	ng	0.00
87) Perylene-d12	15.27	264	305028	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	277664	25.91	ng	-0.01
7) Phenol-d6	6.36	99	378181	28.10	ng	-0.01
23) Nitrobenzene-d5	7.28	82	218885	19.35	ng	-0.02
42) 2,4,6-Tribromophenol	10.55	330	99349	35.86	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	351868	17.48	ng	-0.01
79) Terphenyl-d14	12.82	244	293723	15.93	ng	0.00
Target Compounds						
83) Chrysene	13.89	228	51506	2.122	ng	Qvalue 98
88) Benzo(b)fluoranthene	14.87	252	41857m	2.121	ng	
90) Benzo(a)pyrene	15.22	252	35697	2.045	ng	# 96
92) Benzo(g,h,i)perylene	17.05	276	36239	2.423	ng	91

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

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