

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032219\
 Data File : BF113239.D
 Acq On : 22 Mar 2019 17:22
 Operator : JU/SJ
 Sample : K2004-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-20(10-11)

Manual Integrations
 APPROVED

Sohil
 3/25/2019 6:32:09 PM

Quant Time: Mar 23 00:12:23 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	168414	20.00	ng	-0.02
21) Naphthalene-d8	8.00	136	584978	20.00	ng	-0.02
39) Acenaphthene-d10	9.75	164	286099	20.00	ng	-0.02
64) Phenanthrene-d10	11.23	188	477377	20.00	ng	-0.02
76) Chrysene-d12	13.86	240	346820	20.00	ng	0.00
87) Perylene-d12	15.29	264	399972	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	1129634	104.34	ng	-0.01
7) Phenol-d6	6.36	99	1550019	113.97	ng	-0.01
23) Nitrobenzene-d5	7.28	82	924467	94.56	ng	-0.02
42) 2,4,6-Tribromophenol	10.54	330	428220	140.99	ng	-0.01
45) 2-Fluorobiphenyl	9.07	172	1635670	99.00	ng	-0.02
79) Terphenyl-d14	12.82	244	1292638	72.25	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.46	163	145234	6.885	ng	99
75) Fluoranthene	12.44	202	52772	2.074	ng	94
78) Pyrene	12.66	202	144344	4.937	ng	95
81) Benzo(a)anthracene	13.86	228	96179	4.001	ng	# 80
83) Chrysene	13.89	228	129474	5.499	ng	# 78
88) Benzo(b)fluoranthene	14.88	252	62089m	2.400	ng	
90) Benzo(a)pyrene	15.22	252	83716	3.658	ng	# 61
92) Benzo(g,h,i)perylene	17.06	276	80123	4.086	ng	# 89

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

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