

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012419\
 Data File : BF112213.D
 Acq On : 24 Jan 2019 17:14
 Operator : JU/SJ
 Sample : K1223-01 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-M

Manual Integrations
 APPROVED

Sohil
 1/25/2019 10:05:09 AM

Quant Time: Jan 25 00:23:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 19 01:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	229253	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	893067	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	362699	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	625410	20.00	ng	0.00
76) Chrysene-d12	14.02	240	476903	20.00	ng	0.00
87) Perylene-d12	15.50	264	352245	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	619068	49.01	ng	0.00
7) Phenol-d6	6.50	99	777852	49.22	ng	0.00
23) Nitrobenzene-d5	7.41	82	502054	38.84	ng	-0.02
42) 2,4,6-Tribromophenol	10.68	330	190663	48.73	ng	-0.01
45) 2-Fluorobiphenyl	9.21	172	843192	34.08	ng	0.00
79) Terphenyl-d14	12.96	244	602186	26.86	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.60	163	63903	2.416	ng	97
71) Phenanthrene	11.40	178	170551	5.419	ng	98
75) Fluoranthene	12.59	202	244088	7.345	ng	97
78) Pyrene	12.82	202	215979	6.864	ng	98
81) Benzo(a)anthracene	14.02	228	104463	3.820	ng	95
83) Chrysene	14.05	228	102475	3.961	ng	97
88) Benzo(b)fluoranthene	15.07	252	100842m	5.043	ng	
90) Benzo(a)pyrene	15.44	252	65793	3.588	ng	# 81
92) Benzo(g,h,i)perylene	17.44	276	35927	2.239	ng	# 87

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

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