

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012419\
 Data File : BF112220.D
 Acq On : 24 Jan 2019 20:29
 Operator : JU/SJ
 Sample : K1223-08 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-T

Manual Integrations
 APPROVED

Sohil
 1/25/2019 10:05:23 AM

Quant Time: Jan 25 00:48:05 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	6.86	152	159746	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	636257	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	364230	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	596384	20.00	ng	0.00
76) Chrysene-d12	14.03	240	404210	20.00	ng	0.00
87) Perylene-d12	15.51	264	286912	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	417375	47.42	ng	0.00
7) Phenol-d6	6.50	99	522597	47.46	ng	0.00
23) Nitrobenzene-d5	7.42	82	318262	34.56	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	173852	44.25	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	793575	31.94	ng	0.00
79) Terphenyl-d14	12.96	244	595998	31.37	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.60	163	56692	2.134	ng	99
71) Phenanthrene	11.41	178	145154	4.836	ng	94
75) Fluoranthene	12.60	202	202417	6.388	ng	99
78) Pyrene	12.83	202	172167	6.456	ng	97
81) Benzo(a)anthracene	14.02	228	71462	3.083	ng	95
83) Chrysene	14.06	228	70977	3.237	ng	95
88) Benzo(b)fluoranthene	15.08	252	72557m	4.455	ng	
90) Benzo(a)pyrene	15.44	252	43257	2.896	ng	# 73
92) Benzo(g,h,i)perylene	17.46	276	26580	2.034	ng	# 85

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

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