

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122619\
 Data File : BF118332.D
 Acq On : 26 Dec 2019 22:36
 Operator : JU
 Sample : K6436-07 2X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 K009-WC-2

Manual Integrations
 APPROVED

mohammad
 12/27/2019 3:19:39 PM

Quant Time: Dec 27 06:12:20 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 15:58:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	74220	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	300281	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	152833	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	249858	20.00	ng	0.00
76) Chrysene-d12	14.05	240	190223	20.00	ng	0.00
87) Perylene-d12	15.54	264	147668	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	302789	69.64	ng	0.00
7) Phenol-d6	6.48	99	402861	74.69	ng	0.00
23) Nitrobenzene-d5	7.42	82	226539	43.15	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	88449	46.86	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	516513	51.59	ng	0.00
79) Terphenyl-d14	12.98	244	528706	46.11	ng	0.00
Target Compounds						
10) Phenol	6.50	94	12211	2.014	ng	83
50) Dimethylphthalate	9.61	163	26047	2.270	ng	# 99
71) Phenanthrene	11.42	178	31070	2.281	ng	97
75) Fluoranthene	12.62	202	110804	8.060	ng	99
78) Pyrene	12.84	202	94507	6.084	ng	99
81) Benzo(a)anthracene	14.04	228	42861	3.219	ng	# 96
83) Chrysene	14.07	228	50456	3.958	ng	# 92
88) Benzo(b)fluoranthene	15.10	252	45137m	4.605	ng	
90) Benzo(a)pyrene	15.48	252	27501	3.185	ng	# 41
92) Benzo(g,h,i)perylene	17.52	276	18264	2.523	ng	# 68

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

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