

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012420\
 Data File : BF118688.D
 Acq On : 24 Jan 2020 16:24
 Operator : CG/JU
 Sample : L1237-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PATERSON-UST

Manual Integrations
 APPROVED

mohammad
 1/28/2020 8:12:19 AM

Quant Time: Jan 27 01:40:46 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 20 14:45:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	278508	20.00	ng	-0.02
21) Naphthalene-d8	8.15	136	981848	20.00	ng	-0.02
39) Acenaphthene-d10	9.91	164	414465	20.00	ng	-0.02
64) Phenanthrene-d10	11.40	188	743565	20.00	ng	-0.01
76) Chrysene-d12	14.02	240	604920	20.00	ng	-0.03
87) Perylene-d12	15.47	264	655949	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	1464198	97.64	ng	-0.02
7) Phenol-d6	6.49	99	1806603	92.69	ng	-0.02
23) Nitrobenzene-d5	7.42	82	1140689	64.98	ng	-0.02
42) 2,4,6-Tribromophenol	10.70	330	506939	105.75	ng	-0.01
45) 2-Fluorobiphenyl	9.22	172	1857098	73.65	ng	-0.02
79) Terphenyl-d14	12.97	244	2175515	78.44	ng	-0.02
Target Compounds						
10) Phenol	6.50	94	53667	2.416	ng	97
37) 2-Methylnaphthalene	8.86	142	85557	2.733	ng	# 92
38) 1-Methylnaphthalene	8.96	142	135167	4.544	ng	# 98
50) Dimethylphthalate	9.62	163	340679	12.616	ng	# 88
52) Acenaphthene	9.94	154	70058	3.093	ng	# 79
71) Phenanthrene	11.42	178	593874	16.555	ng	# 90
78) Pyrene	12.83	202	117182m	2.853	ng	

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

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