

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101419\
 Data File : BF117404.D
 Acq On : 14 Oct 2019 17:53
 Operator : JU
 Sample : PB123882BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB123882BL

Manual Integrations
 APPROVED

mohammad
 10/16/2019 8:10:33 AM

Quant Time: Oct 15 14:25:18 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 14 19:03:29 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	58183	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	236062	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	125403	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	214768	20.00	ng	0.00
76) Chrysene-d12	13.96	240	152752	20.00	ng	0.00
87) Perylene-d12	15.42	264	137650	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	296516	78.41	ng	0.00
7) Phenol-d6	6.43	99	261176	52.06	ng	0.00
23) Nitrobenzene-d5	7.36	82	436099	93.65	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	103934	101.83	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	741906	86.25	ng	0.00
79) Terphenyl-d14	12.90	244	820141	87.53	ng	0.00
Target Compounds						
2) 1,4-Dioxane	2.58	88	5207	3.345	ng	Qvalue # 90
3) Pyridine	3.36	79	27951	7.055	ng	98
4) n-Nitrosodimethylamine	3.30	42	3059m	2.191	ng	
6) Aniline	6.46	93	64300	10.259	ng	# 73
15) Benzyl Alcohol	6.95	79	17484	4.555	ng	# 43
33) 4-Chloroaniline	8.17	127	24201	4.834	ng	98
77) Benzidine	12.65	184	59803	15.830	ng	96

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

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