

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011719\
 Data File : BM018610.D
 Acq On : 17 Jan 2019 12:48
 Operator : JU/SJ
 Sample : K1149-02
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SW002-190114-01

Manual Integrations
 APPROVED

Sohil
 1/18/2019 12:06:52 PM

Quant Time: Jan 18 04:55:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 10 05:57:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	351464	20.00	ng	-0.02
21) Naphthalene-d8	10.53	136	1397212	20.00	ng	-0.02
39) Acenaphthene-d10	14.39	164	741085	20.00	ng	-0.02
64) Phenanthrene-d10	17.14	188	1456320	20.00	ng	-0.02
76) Chrysene-d12	21.33	240	1544012	20.00	ng	-0.02
87) Perylene-d12	23.61	264	1425362	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	1263140	66.36	ng	-0.02
7) Phenol-d6	6.92	99	1325191	54.82	ng	-0.02
23) Nitrobenzene-d5	8.91	82	1974831	81.94	ng	-0.01
42) 2,4,6-Tribromophenol	15.89	330	1153033	122.19	ng	-0.02
45) 2-Fluorobiphenyl	13.01	172	4774227	84.06	ng	-0.01
79) Terphenyl-d14	19.77	244	6212235	84.82	ng	-0.01
Target Compounds						
3) Pyridine	3.68	79	655072	33.869	ng	Qvalue 96
17) 2-Methylphenol	8.19	107	37664	2.259	ng	97
20) 3+4-Methylphenols	8.52	107	53739	2.334	ng	87
80) Butylbenzylphthalate	20.47	149	84801	2.171	ng	94
81) Benzo(a)anthracene	21.32	228	216894	2.384	ng	97
83) Chrysene	21.37	228	212727m	2.422	ng	
84) Bis(2-ethylhexyl)phthalate	21.24	149	118360	2.099	ng	98
88) Benzo(b)fluoranthene	22.92	252	167012	2.064	ng	# 88

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

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