

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121318\
 Data File : BF111383.D
 Acq On : 13 Dec 2018 21:25
 Operator : JU/SJ
 Sample : J6305-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SW003-01

Manual Integrations
 APPROVED

Sohil
 12/14/2018 12:53:36 PM

Quant Time: Dec 14 00:14:33 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 13:16:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	334552	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	1218843	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	500509	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	668151	20.00	ng	0.00
76) Chrysene-d12	13.96	240	479791	20.00	ng	0.00
87) Perylene-d12	15.40	264	341486	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	1220328	61.68	ng	0.02
7) Phenol-d6	6.46	99	805592m	30.20	ng	0.03
23) Nitrobenzene-d5	7.37	82	1787052	82.64	ng	0.00
42) 2,4,6-Tribromophenol	10.64	330	454937	103.11	ng	0.02
45) 2-Fluorobiphenyl	9.16	172	2710967	86.96	ng	0.00
79) Terphenyl-d14	12.91	244	1663475	76.31	ng	0.00
Target Compounds						
2) 1,4-Dioxane	2.53	88	149108	15.089	ng	Qvalue # 91
3) Pyridine	3.29	79	162993	6.714	ng	91
10) Phenol	6.52	94	11231528m	371.706	ng	
17) 2-Methylphenol	7.09	107	1223301	62.257	ng	98
20) 3+4-Methylphenols	7.26	107	1668366	71.198	ng	# 61
22) Acetophenone	7.22	105	791800	27.967	ng	# 85
27) 2,4-Dimethylphenol	7.77	122	498989	30.023	ng	96
31) Naphthalene	8.10	128	296182	5.216	ng	98
84) Bis(2-ethylhexyl)phthalate	13.94	149	136338	6.273	ng	98

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

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