

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121318\
 Data File : BF111388.D
 Acq On : 13 Dec 2018 23:43
 Operator : JU/SJ
 Sample : J6305-08
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SW005-01

Manual Integrations
 APPROVED

Sohil
 12/14/2018 12:53:44 PM

Quant Time: Dec 14 10:55:34 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	232892	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	1161987	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	472725	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	645631	20.00	ng	0.00
76) Chrysene-d12	13.95	240	470081	20.00	ng	0.00
87) Perylene-d12	15.39	264	361172	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	981071	71.24	ng	0.03
7) Phenol-d6	6.46	99	168340	9.07	ng	0.03
23) Nitrobenzene-d5	7.37	82	1457180	70.68	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	395366	94.88	ng	0.01
45) 2-Fluorobiphenyl	9.16	172	2285494	77.62	ng	0.00
79) Terphenyl-d14	12.91	244	1545111	72.34	ng	0.00
Target Compounds						
2) 1,4-Dioxane	2.53	88	387282	56.297	ng	96
3) Pyridine	3.29	79	325449	19.258	ng	100
6) Aniline	6.47	93	3618844	161.097	ng	# 51
10) Phenol	6.53	94	15921966	756.949	ng	99
17) 2-Methylphenol	7.10	107	1995320	145.874	ng	98
20) 3+4-Methylphenols	7.26	107	1916109m	117.465	ng	
22) Acetophenone	7.23	105	1796417	66.555	ng	# 87
27) 2,4-Dimethylphenol	7.76	122	622488	39.287	ng	98
31) Naphthalene	8.10	128	216499	3.999	ng	96
50) Dimethylphthalate	9.56	163	163986	4.824	ng	# 64

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

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