

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121318\
 Data File : BF111384.D
 Acq On : 13 Dec 2018 21:53
 Operator : JU/SJ
 Sample : J6305-04
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SW004-01

Manual Integrations
 APPROVED

Sohil
 12/17/2018 2:10:14 PM

Quant Time: Dec 17 10:38:32 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	308379	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	1271013	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	490669	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	688720	20.00	ng	0.00
76) Chrysene-d12	13.95	240	478114	20.00	ng	0.00
87) Perylene-d12	15.39	264	382750	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	901476	49.43	ng	0.02
7) Phenol-d6	6.47	99	658218	26.77	ng	0.04
23) Nitrobenzene-d5	7.36	82	1383390	61.35	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	339540	78.50	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	2003397	65.55	ng	0.00
79) Terphenyl-d14	12.90	244	1091702	50.26	ng	0.00
Target Compounds						
2) 1,4-Dioxane	2.56	88	29184	3.204	ng	# 79
10) Phenol	6.49	94	5403288	193.998	ng	97
17) 2-Methylphenol	7.08	107	569318	31.433	ng	98
20) 3+4-Methylphenols	7.24	107	810082	37.505	ng	# 59
22) Acetophenone	7.20	105	186745	6.325	ng	# 87
27) 2,4-Dimethylphenol	7.76	122	132741	7.659	ng	# 92
29) 2,4-Dichlorophenol	7.96	162	106753	5.905	ng	98
32) Benzoic acid	8.02	122	2083678m	144.625	ng	
35) Caprolactam	8.65	113	1812140m	286.052	ng	
50) Dimethylphthalate	9.56	163	153824	4.360	ng	# 88

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

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