

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121318\
 Data File : BF111387.D
 Acq On : 13 Dec 2018 23:16
 Operator : JU/SJ
 Sample : J6305-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SW004-02

Manual Integrations
 APPROVED

Sohil
 12/17/2018 2:10:16 PM

Quant Time: Dec 17 10:44:56 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	338713	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	1345633	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	557166	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	753696	20.00	ng	0.00
76) Chrysene-d12	13.95	240	539377	20.00	ng	0.00
87) Perylene-d12	15.39	264	446637	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	752346	37.56	ng	0.02
7) Phenol-d6	6.47	99	526615	19.50	ng	0.04
23) Nitrobenzene-d5	7.36	82	1114187	46.67	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	290816	59.21	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1756582	50.62	ng	0.00
79) Terphenyl-d14	12.90	244	1010109	41.22	ng	0.00
Target Compounds						
2) 1,4-Dioxane	2.55	88	26570	2.656	ng	# 55
10) Phenol	6.49	94	5098831	166.672	ng	96
17) 2-Methylphenol	7.08	107	463871	23.318	ng	97
20) 3+4-Methylphenols	7.24	107	649506	27.378	ng	# 59
22) Acetophenone	7.20	105	164201	5.253	ng	# 85
29) 2,4-Dichlorophenol	7.96	162	72877	3.807	ng	97
32) Benzoic acid	8.06	122	988114m	64.781	ng	
35) Caprolactam	8.65	113	1981049m	295.373	ng	
50) Dimethylphthalate	9.56	163	150566	3.758	ng	# 84

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

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