

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121418\
 Data File : BF111447.D
 Acq On : 15 Dec 2018 4:49
 Operator : JU/SJ
 Sample : J6305-08DL 20X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SW005-01DL

Manual Integrations
 APPROVED

Sohil
 12/17/2018 2:37:19 PM

Quant Time: Dec 17 10:51:06 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	138084	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	521437	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	209908	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	323888	20.00	ng	0.00
76) Chrysene-d12	13.95	240	266462	20.00	ng	0.00
87) Perylene-d12	15.39	264	183228	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	23669	2.85	ng	0.00
7) Phenol-d6	6.43	99	18000	1.63	ng	0.00
23) Nitrobenzene-d5	7.35	82	39905	4.56	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	9549	5.08	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	74810	5.51	ng	0.00
79) Terphenyl-d14	12.90	244	53586	4.72	ng	0.00
Target Compounds						
2) 1,4-Dioxane	2.48	88	11470	2.862	ng	89
6) Aniline	6.45	93	138472m	10.114	ng	
10) Phenol	6.45	94	567286	46.136	ng	98
17) 2-Methylphenol	7.05	107	58436	7.323	ng	97
20) 3+4-Methylphenols	7.21	107	51475	5.143	ng	# 64
22) Acetophenone	7.20	105	50283	4.320	ng	# 90
27) 2,4-Dimethylphenol	7.73	122	16371	2.438	ng	96

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

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