Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121418\
Data File : BF111449.D
Acq On : 15 Dec 2018 5:44
Operator : JU/SJ
Sample : J6305-03DL 10X
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : P001-SW003-01DL

Quant Time: Dec 15 23:47:30 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	148240	20.00	ng		0.00
21) Naphthalene-d8	8.07	136	549423	20.00	ng		0.00
39) Acenaphthene-d10	9.83	164	215576	20.00	ng		0.00
64) Phenanthrene-d10	11.32	188	342882	20.00	ng		0.00
76) Chrysene-d12	13.95	240	287147	20.00	ng		0.00
87) Perylene-d12	15.39	264	190503	20.00	ng		0.00
System Monitoring Compounds							
5) 2-Fluorophenol	5.42	112	66256	7.44	ng		0.00
7) Phenol-d6	6.43	99	49715	4.20	ng		0.00
23) Nitrobenzene-d5	7.35	82	99118	10.74	ng		0.00
42) 2,4,6-Tribromophenol	10.62	330	26596	13.77	ng		0.00
45) 2-Fluorobiphenyl	9.15	172	182595	13.08	ng		0.00
79) Terphenyl-d14	12.90	244	132379	10.83	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	2.49	88	9137	2.124	ng	#	86
10) Phenol	6.45	94	762932	57.797	ng		98
17) 2-Methylphenol	7.06	107	67981	7.936	ng		98
20) 3+4-Methylphenols	7.21	107	107291	9.985	ng	#	60
22) Acetophenone	7.20	105	45143	3.681	ng	#	90
27) 2,4-Dimethylphenol	7.73	122	24308	3.435	ng		95

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

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