

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
 Data File : BF111726.D
 Acq On : 26 Dec 2018 21:03
 Operator : JU/SJ
 Sample : J6501-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VNJ-221

Manual Integrations
 APPROVED

Sohil
 12/27/2018 8:22:31 AM

Quant Time: Dec 27 02:56:57 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	210077	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	763589	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	344036	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	566027	20.00	ng	0.00
76) Chrysene-d12	14.06	240	501304	20.00	ng	0.00
87) Perylene-d12	15.54	264	378775	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	751362	60.96	ng	0.02
7) Phenol-d6	6.53	99	946555	60.35	ng	0.01
23) Nitrobenzene-d5	7.46	82	587299	44.77	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	228370	56.18	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	1002294	42.46	ng	0.00
79) Terphenyl-d14	13.00	244	874023	33.06	ng	0.00
Target Compounds						
10) Phenol	6.55	94	70123	3.962	ng	98
50) Dimethylphthalate	9.65	163	124848	4.962	ng	97
75) Fluoranthene	12.63	202	152584	5.020	ng	92
78) Pyrene	12.86	202	138251	3.905	ng	95
81) Benzo(a)anthracene	14.04	228	77145	2.697	ng	96
83) Chrysene	14.08	228	80056	2.847	ng	97
88) Benzo(b)fluoranthene	15.10	252	99722m	4.505	ng	
90) Benzo(a)pyrene	15.47	252	63324	3.149	ng	95
92) Benzo(g,h,i)perylene	17.47	276	39713	2.309	ng	# 86

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

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