

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102618\
 Data File : BF110211.D
 Acq On : 26 Oct 2018 15:56
 Operator : JU/SJ
 Sample : J5621-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SA-SB-01A

Manual Integrations
 APPROVED

Sohil
 10/29/2018 11:56:39 AM

Quant Time: Oct 26 23:55:18 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 23 15:06:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	122447	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	480401	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	198799	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	310258	20.00	ng	-0.01
76) Chrysene-d12	14.15	240	226223	20.00	ng	-0.01
87) Perylene-d12	15.68	264	160128	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	748590	99.56	ng	0.00
7) Phenol-d6	6.59	99	969611	100.82	ng	-0.01
23) Nitrobenzene-d5	7.53	82	600287	74.12	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	167829	85.23	ng	-0.01
45) 2-Fluorobiphenyl	9.33	172	928922	73.37	ng	-0.01
79) Terphenyl-d14	13.09	244	641235	58.95	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	136604	13.288	ng	Qvalue # 64
50) Dimethylphthalate	9.72	163	151162	10.301	ng	99
71) Phenanthrene	11.53	178	36701	2.261	ng	99
75) Fluoranthene	12.72	202	61624	3.740	ng	97
78) Pyrene	12.95	202	58524	3.609	ng	98
81) Benzo(a)anthracene	14.14	228	27862	2.077	ng	91
83) Chrysene	14.17	228	37427	2.937	ng	96
88) Benzo(b)fluoranthene	15.22	252	38518m	4.166	ng	
90) Benzo(a)pyrene	15.61	252	19619	2.306	ng	# 85

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

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