

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102518\
 Data File : BF110192.D
 Acq On : 26 Oct 2018 6:16
 Operator : JU/SJ
 Sample : J5658-02
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-TS-002

Manual Integrations
 APPROVED

Sohil
 10/26/2018 3:49:35 PM

Quant Time: Oct 26 09:23:49 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	141377	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	531501	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	214510	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	349098	20.00	ng	-0.01
76) Chrysene-d12	14.15	240	242793	20.00	ng	-0.01
87) Perylene-d12	15.68	264	189382	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	851976	98.13	ng	0.00
7) Phenol-d6	6.59	99	1059610	95.42	ng	-0.01
23) Nitrobenzene-d5	7.53	82	626934	69.96	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	182478	85.88	ng	-0.01
45) 2-Fluorobiphenyl	9.33	172	908800	66.53	ng	-0.01
79) Terphenyl-d14	13.09	244	627938	53.79	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	119779	10.091	ng	Qvalue # 64
50) Dimethylphthalate	9.72	163	157445	9.943	ng	100
75) Fluoranthene	12.72	202	66225	3.572	ng	97
78) Pyrene	12.95	202	55266	3.175	ng	99
83) Chrysene	14.17	228	31123	2.276	ng	97
88) Benzo(b)fluoranthene	15.22	252	40387m	3.693	ng	
90) Benzo(a)pyrene	15.61	252	25552	2.539	ng	# 92

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

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