

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102618\
 Data File : BF110226.D
 Acq On : 26 Oct 2018 22:45
 Operator : JU/SJ
 Sample : J5675-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-13

Manual Integrations
 APPROVED

Sohil
 10/29/2018 11:56:48 AM

Quant Time: Oct 27 00:46:47 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	124427	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	484425	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	190650	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	300216	20.00	ng	-0.01
76) Chrysene-d12	14.15	240	212617	20.00	ng	-0.01
87) Perylene-d12	15.68	264	146826	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	734864	96.18	ng	0.00
7) Phenol-d6	6.59	99	915736	93.70	ng	-0.01
23) Nitrobenzene-d5	7.53	82	535296	65.54	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	148907	78.85	ng	-0.01
45) 2-Fluorobiphenyl	9.33	172	841845	69.34	ng	-0.01
79) Terphenyl-d14	13.09	244	619742	60.62	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	64612	6.185	ng	Qvalue # 64
50) Dimethylphthalate	9.72	163	151293	10.750	ng	99
75) Fluoranthene	12.72	202	79928	5.014	ng	99
78) Pyrene	12.95	202	79483	5.214	ng	97
81) Benzo(a)anthracene	14.14	228	44069	3.495	ng	97
83) Chrysene	14.17	228	41052	3.428	ng	97
86) Indeno(1,2,3-cd)pyrene	17.24	276	23038	2.319	ng	94
88) Benzo(b)fluoranthene	15.22	252	48901m	5.768	ng	
90) Benzo(a)pyrene	15.62	252	31940	4.094	ng	# 94
92) Benzo(g,h,i)perylene	17.72	276	22214	3.349	ng	# 86

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

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