

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110918\
 Data File : BF110645.D
 Acq On : 9 Nov 2018 23:25
 Operator : JU/SJ
 Sample : J5876-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 48382-47191-47202

Manual Integrations
 APPROVED

Sohil
 11/12/2018 10:30:24 AM

Quant Time: Nov 11 03:58:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 09 15:13:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	88689	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	349671	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	147132	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	223386	20.00	ng	0.00
76) Chrysene-d12	14.13	240	159746	20.00	ng	0.00
87) Perylene-d12	15.66	264	125112	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	731004	133.88	ng	0.01
7) Phenol-d6	6.58	99	941676	134.87	ng	0.00
23) Nitrobenzene-d5	7.52	82	594856	97.97	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	177755	119.90	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	918498	89.16	ng	0.00
79) Terphenyl-d14	13.06	244	592578	69.47	ng	0.00
Target Compounds						
10) Phenol	6.59	94	58161	7.184	ng	Qvalue # 69
50) Dimethylphthalate	9.69	163	56974	5.190	ng	100
71) Phenanthrene	11.50	178	94815	7.922	ng	99
75) Fluoranthene	12.70	202	128124	10.678	ng	98
78) Pyrene	12.93	202	118272	9.616	ng	96
81) Benzo(a)anthracene	14.12	228	57840	6.058	ng	96
83) Chrysene	14.16	228	53173	5.845	ng	98
84) Bis(2-ethylhexyl)phthalate	14.08	149	23108	3.519	ng	# 92
86) Indeno(1,2,3-cd)pyrene	17.23	276	25598	3.921	ng	94
88) Benzo(b)fluoranthene	15.20	252	65061m	8.693	ng	
89) Benzo(k)fluoranthene	15.23	252	23677m	3.201	ng	
90) Benzo(a)pyrene	15.59	252	44740	6.579	ng	# 92
92) Benzo(a,h,i)perylene	17.70	276	23460	4.109	ng	# 87

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

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