

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122018\
 Data File : BF111643.D
 Acq On : 20 Dec 2018 17:19
 Operator : JU/SJ
 Sample : J6445-06
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EB05_1.5-3.5

Manual Integrations
 APPROVED

Sohil
 12/21/2018 2:46:43 PM

Quant Time: Dec 21 06:42:10 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	210235	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	794690	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	405168	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	729587	20.00	ng	0.00
76) Chrysene-d12	14.05	240	460849	20.00	ng	0.00
87) Perylene-d12	15.53	264	489433	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	907297	73.56	ng	0.00
7) Phenol-d6	6.52	99	1187832	75.67	ng	0.00
23) Nitrobenzene-d5	7.45	82	732260	53.63	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	295255	61.67	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1479498	53.22	ng	0.00
79) Terphenyl-d14	13.00	244	1283282	52.81	ng	0.00
Target Compounds						
10) Phenol	6.53	94	74631	4.214	ng	91
50) Dimethylphthalate	9.64	163	107225	3.619	ng	98
71) Phenanthrene	11.44	178	688057	17.783	ng	94
72) Anthracene	11.49	178	123669	3.184	ng	93
75) Fluoranthene	12.63	202	1011629	25.819	ng	95
78) Pyrene	12.85	202	885414	27.207	ng	96
81) Benzo(a)anthracene	14.04	228	428133	16.279	ng	99
83) Chrysene	14.07	228	397692	15.385	ng	98
86) Indeno(1,2,3-cd)pyrene	17.00	276	254925	11.601	ng	# 82
88) Benzo(b)fluoranthene	15.10	252	599960m	20.974	ng	
89) Benzo(k)fluoranthene	15.12	252	150318m	5.520	ng	
90) Benzo(a)pyrene	15.46	252	396884m	15.272	ng	
91) Dibenzo(a,h)anthracene	17.02	278	64075	2.816	ng	# 91
92) Benzo(a,h,i)perylene	17.45	276	219222	9.866	ng	# 88

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

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