

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
 Data File : BF111722.D
 Acq On : 26 Dec 2018 19:14
 Operator : JU/SJ
 Sample : J6499-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 72-11989

Manual Integrations
 APPROVED

Sohil
 12/27/2018 8:22:24 AM

Quant Time: Dec 27 02:39:03 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	247004	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	941016	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	482774	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	804414	20.00	ng	0.00
76) Chrysene-d12	14.06	240	541003	20.00	ng	0.00
87) Perylene-d12	15.54	264	536730	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	1431706	98.80	ng	0.02
7) Phenol-d6	6.54	99	1921831	104.21	ng	0.02
23) Nitrobenzene-d5	7.46	82	1256382	77.72	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	554470	97.20	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	2059031	62.17	ng	0.00
79) Terphenyl-d14	13.00	244	1681817	58.95	ng	0.00
Target Compounds						
10) Phenol	6.55	94	162342	7.802	ng	99
50) Dimethylphthalate	9.65	163	312509	8.852	ng	97
71) Phenanthrene	11.45	178	128269	3.007	ng	93
75) Fluoranthene	12.63	202	362722	8.396	ng	94
78) Pyrene	12.86	202	322568	8.443	ng	96
81) Benzo(a)anthracene	14.04	228	190807	6.180	ng	99
83) Chrysene	14.08	228	166609	5.490	ng	98
86) Indeno(1,2,3-cd)pyrene	17.02	276	62918	2.439	ng	# 85
88) Benzo(b)fluoranthene	15.10	252	214101m	6.825	ng	
90) Benzo(a)pyrene	15.47	252	140873m	4.943	ng	
92) Benzo(g,h,i)perylene	17.47	276	53242m	2.185	ng	

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

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