

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
 Data File : BF111995.D
 Acq On : 10 Jan 2019 18:49
 Operator : JU/SJ
 Sample : K1071-04
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CPSB-16(20-25)

Manual Integrations
 APPROVED

Sohil
 1/11/2019 4:42:53 PM

Quant Time: Jan 11 01:36:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 07 14:33:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	102228	20.00	ng	-0.02
21) Naphthalene-d8	8.16	136	327556	20.00	ng	-0.01
39) Acenaphthene-d10	9.92	164	143039	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	120368	20.00	ng	-0.02
76) Chrysene-d12	14.04	240	133683	20.00	ng	-0.02
87) Perylene-d12	15.53	264	107930	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	578810	96.87	ng	-0.01
7) Phenol-d6	6.52	99	743138	96.49	ng	-0.02
23) Nitrobenzene-d5	7.43	82	504099	82.25	ng	-0.02
42) 2,4,6-Tribromophenol	10.70	330	81964	44.05	ng	-0.02
45) 2-Fluorobiphenyl	9.24	172	716277	73.02	ng	-0.01
79) Terphenyl-d14	12.98	244	346060	49.14	ng	-0.02
Target Compounds						
10) Phenol	6.53	94	89281	10.666	ng	# 77
20) 3+4-Methylphenols	7.28	107	112997	17.026	ng	# 65
50) Dimethylphthalate	9.63	163	99233	9.235	ng	# 86
52) Acenaphthene	9.95	154	106108	11.855	ng	95
55) Dibenzofuran	10.12	168	112952	8.678	ng	# 90
58) Fluorene	10.46	166	60501	6.388	ng	97
71) Phenanthrene	11.42	178	182385	28.341	ng	98
72) Anthracene	11.47	178	20831	3.189	ng	98
75) Fluoranthene	12.61	202	83925	12.703	ng	99
78) Pyrene	12.84	202	64970	7.063	ng	99
81) Benzo(a)anthracene	14.03	228	18750	2.437	ng	100
83) Chrysene	14.06	228	15172m	2.068	ng	
84) Bis(2-ethylhexyl)phthalate	14.02	149	377271	73.873	ng	99

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

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