

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011119\
 Data File : BF112032.D
 Acq On : 11 Jan 2019 22:55
 Operator : JU/SJ
 Sample : K1102-03 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CPSB-02(15-20)

Manual Integrations
 APPROVED

Sohil
 1/14/2019 4:10:42 PM

Quant Time: Jan 12 01:55:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 11 21:37:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	17771	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	59672	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	27863	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	45008	20.00	ng	0.01
76) Chrysene-d12	14.07	240	30280	20.00	ng	0.02
87) Perylene-d12	15.56	264	30458	20.00	ng	0.04

System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	107234	103.23	ng	0.00
7) Phenol-d6	6.52	99	136337	101.83	ng	0.00
23) Nitrobenzene-d5	7.43	82	77824	69.70	ng	-0.01
42) 2,4,6-Tribromophenol	10.71	330	30968	85.45	ng	0.00
45) 2-Fluorobiphenyl	9.23	172	144863	75.81	ng	0.00
79) Terphenyl-d14	13.00	244	130064	81.54	ng	0.02

Target Compounds				Qvalue		
38) 1-Methylnaphthalene	8.97	142	9935	5.901	ng	# 91
50) Dimethylphthalate	9.62	163	37935	18.124	ng	# 87
52) Acenaphthene	9.95	154	129685	74.385	ng	# 98
55) Dibenzofuran	10.12	168	79868	31.500	ng	# 94
58) Fluorene	10.46	166	116446	63.116	ng	# 94
71) Phenanthrene	11.43	178	111597	46.376	ng	# 89
72) Anthracene	11.49	178	58613	23.994	ng	# 82
75) Fluoranthene	12.63	202	158894	64.321	ng	# 91
78) Pyrene	12.87	202	132597m	63.643	ng	# 91
81) Benzo(a)anthracene	14.06	228	32463	18.628	ng	# 64
83) Chrysene	14.09	228	32324	19.454	ng	# 66
84) Bis(2-ethylhexyl)phthalate	14.03	149	71942	62.192	ng	# 78
86) Indeno(1,2,3-cd)pyrene	17.06	276	10398	8.584	ng	# 85
88) Benzo(b)fluoranthene	15.12	252	23090m	12.673	ng	# 85
89) Benzo(k)fluoranthene	15.14	252	9385m	5.480	ng	# 85
90) Benzo(a)pyrene	15.49	252	14855	9.273	ng	# 1
92) Benzo(a,h,i)perylene	17.51	276	10115	8.026	ng	# 4

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

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