

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042219\
 Data File : BF113906.D
 Acq On : 23 Apr 2019 3:58
 Operator : JU/SJ
 Sample : K2201-09RE 5X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-01-041819-E

Quant Time: Apr 23 06:41:34 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF042219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 23 03:32:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	0.00	152	0	0.00	ng	-6.90
21) Naphthalene-d8	0.00	136	0	0.00	ng	-8.18
39) Acenaphthene-d10	0.00	164	0	0.00	ng	-9.93
64) Phenanthrene-d10	0.00	188	0	0.00	ng	-11.42
76) Chrysene-d12	14.06	240	1739	20.00	ng	-0.01
87) Perylene-d12	15.53	264	452	20.00	ng	-0.04

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	0.00	112	0	0.00	ng	
7) Phenol-d6	0.00	99	0	0.00	ng	
23) Nitrobenzene-d5	0.00	82	0	0.00	ng	
42) 2,4,6-Tribromophenol	0.00	330	0	0.00	ng	
45) 2-Fluorobiphenyl	0.00	172	0	0.00	ng	
79) Terphenyl-d14	13.00	244	1099	12.96	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
78) Pyrene	12.86	202	546	4.490	ng	# 10
80) Butylbenzylphthalate	13.49	149	233	4.870	ng	# 66
81) Benzo(a)anthracene	14.04	228	1838	17.939	ng	# 1
82) 3,3'-Dichlorobenzidine	13.99	252	1151	30.649	ng	# 25
83) Chrysene	14.08	228	2359	23.641	ng	# 56
84) Bis(2-ethylhexyl)phthalate	14.02	149	235	3.860	ng	# 51
85) Di-n-octyl phthalate	14.72	149	944	9.907	ng	# 62
86) Indeno(1,2,3-cd)pyrene	17.02	276	277	2.931	ng	# 1
88) Benzo(b)fluoranthene	15.10	252	755	26.401	ng	# 1
89) Benzo(k)fluoranthene	15.13	252	837	34.026	ng	# 1
90) Benzo(a)pyrene	15.47	252	448	18.111	ng	# 1
91) Dibenzo(a,h)anthracene	17.04	278	317	13.652	ng	# 1
92) Benzo(g,h,i)perylene	17.46	276	535	22.831	ng	# 1

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

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