

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090119\
 Data File : BG042665.D
 Acq On : 1 Sep 2019 16:03
 Operator : HP/JU
 Sample : DFTPP
 Misc :
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DFTPP

Quant Time: Sep 03 12:16:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 22 14:29:15 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	-8.01
21) Naphthalene-d8	0.00	136	0	0.00	ng	-10.83
39) Acenaphthene-d10	0.00	164	0	0.00	ng	-14.66
64) Phenanthrene-d10	17.42	188	217	20.00	ng	0.00
76) Chrysene-d12	21.70	240	526	20.00	ng	0.00
87) Perylene-d12	24.95	264	508	20.00	ng	0.01

System Monitoring Compounds

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	20.02	244	1259	51.75	ng	0.00

Target Compounds

						Qvalue
69) Atrazine	17.07	200	19641	9306.997	ng	# 35
70) Pentachlorophenol	17.07	266	128646	82747.267	ng	# 99
71) Phenanthrene	17.46	178	320	29.688	ng	# 59
72) Anthracene	17.56	178	182	16.914	ng	# 60
73) Carbazole	17.85	167	6427	670.167	ng	# 85
74) Di-n-butylphthalate	18.38	149	469	41.371	ng	# 77
75) Fluoranthene	19.47	202	532	40.442	ng	# 63
77) Benzidine	19.64	184	753156	49941.902	ng	# 99
78) Pyrene	19.83	202	666	20.652	ng	# 55
80) Butylbenzylphthalate	20.71	149	271	21.365	ng	# 70
81) Benzo(a)anthracene	21.68	228	554	17.314	ng	# 49
82) 3,3'-Dichlorobenzidine	21.60	252	801	62.243	ng	# 90
83) Chrysene	21.75	228	749	24.272	ng	# 48
84) Bis(2-ethylhexyl)phthalate	21.56	149	1193	67.741	ng	# 87
85) Di-n-octyl phthalate	22.78	149	768	25.355	ng	# 84
86) Indeno(1,2,3-cd)pyrene	28.70	276	182	4.340	ng	# 55
88) Benzo(b)fluoranthene	23.91	252	251	8.987	ng	# 59
89) Benzo(k)fluoranthene	24.00	252	946	35.240	ng	# 1
90) Benzo(a)pyrene	24.79	252	61	2.302	ng	# 59
92) Benzo(a,h,i)perylene	29.83	276	61	2.273	ng	# 56

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

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