Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092619\
Data File : BG042936.D
Acq On : 26 Sep 2019 13:27
Operator : JU
Sample : K4839-06DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : PT-ACIDS-WPDL

Quant Time: Sep 26 14:24:35 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 25 15:35:52 2019
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	48859	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	186188	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	137098	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	362119	20.00	ng	0.00
76) Chrysene-d12	21.71	240	407207	20.00	ng	0.00
87) Perylene-d12	24.94	264	447189	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.65	112	86511	31.60	ng	0.00
7) Phenol-d6	7.24	99	127984	32.46	ng	0.00
23) Nitrobenzene-d5	9.23	82	82454	21.52	ng	0.00
42) 2,4,6-Tribromophenol	16.18	330	76040	32.25	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	226713	23.97	ng	0.00
79) Terphenyl-d14	20.04	244	526058	27.53	ng	0.00
Target Compounds						Qvalue
8) 2-Chlorophenol	7.64	128	89887	31.482	ng	89
10) Phenol	7.26	94	129606	35.831	ng	94
20) 3+4-Methylphenols	8.84	107	72815	20.993	ng	93
26) 2-Nitrophenol	9.99	139	55930	35.958	ng	94
27) 2,4-Dimethylphenol	10.05	122	106336	44.516	ng	97
29) 2,4-Dichlorophenol	10.53	162	103911	34.977	ng	98
36) 4-Chloro-3-methylphenol	12.15	107	81924	25.362	ng	99
43) 2,4,6-Trichlorophenol	13.13	196	38894	12.891	ng	97
44) 2,4,5-Trichlorophenol	13.20	196	33846	10.890	ng	95
54) 2,4-Dinitrophenol	14.79	184	40387	29.702	ng	97
56) 4-Nitrophenol	14.91	139	40304	23.398	ng	95
65) 4,6-Dinitro-2-methylphenol	15.81	198	56758	27.699	ng	96
70) Pentachlorophenol	17.09	266	54532	18.675	ng	99

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

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