

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092619\
 Data File : BG042935.D
 Acq On : 26 Sep 2019 12:07
 Operator : JU
 Sample : K4839-06
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PT-ACIDS-WP

Quant Time: Sep 26 14:19:54 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	8.08	152	61978	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	222540	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	165568	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	397715	20.00	ng	0.00
76) Chrysene-d12	21.71	240	459939	20.00	ng	0.00
87) Perylene-d12	24.94	264	496176	20.00	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.66	112	502125	144.58	ng	0.00
7) Phenol-d6	7.25	99	742833	148.53	ng	0.00
23) Nitrobenzene-d5	9.23	82	467045	102.01	ng	0.00
42) 2,4,6-Tribromophenol	16.18	330	432717	151.99	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	1134311	99.32	ng	0.00
79) Terphenyl-d14	20.04	244	2329331	107.92	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
8) 2-Chlorophenol	7.64	128	507619	140.158	ng	98
10) Phenol	7.28	94	721524	157.249	ng	98
20) 3+4-Methylphenols	8.84	107	447774	101.772	ng	92
26) 2-Nitrophenol	9.99	139	342627	184.295	ng	96
27) 2,4-Dimethylphenol	10.06	122	592686	207.589	ng	99
29) 2,4-Dichlorophenol	10.53	162	618117	174.076	ng	97
36) 4-Chloro-3-methylphenol	12.15	107	478904	124.043	ng	98
43) 2,4,6-Trichlorophenol	13.13	196	235678	64.682	ng	99
44) 2,4,5-Trichlorophenol	13.20	196	206828	55.102	ng	98
54) 2,4-Dinitrophenol	14.80	184	304558	185.468	ng	93
56) 4-Nitrophenol	14.91	139	248016	119.226	ng	96
65) 4,6-Dinitro-2-methylphenol	15.81	198	372220	165.392	ng	96
70) Pentachlorophenol	17.09	266	347327	108.301	ng	98

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

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