

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100515\
 Data File : BF082047.D
 Acq On : 5 Oct 2015 15:53
 Operator : UM/IZ
 Sample : G3692-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PT-ACIDS-WP

Manual Integrations
 APPROVED

MMDadoda
 10/6/2015 2:55:11 PM

Quant Time: Oct 05 16:32:03 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 05 15:47:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	107555	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	437589	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	195721	20.00	ng	-0.01
63) Phenanthrene-d10	11.42	188	400271	20.00	ng	0.00
75) Chrysene-d12	14.06	240	375687	20.00	ng	0.00
86) Perylene-d12	15.50	264	319745	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	715547	120.77	ng	0.01
7) Phenol-d6	6.53	99	991381	132.97	ng	0.01
23) Nitrobenzene-d5	7.45	82	592861	90.31	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	296011	149.34	ng	0.00
44) 2-Fluorobiphenyl	9.25	172	1166257	104.24	ng	0.00
78) Terphenyl-d14	13.01	244	1277130	126.14	ng	0.00
Target Compounds						Qvalue
8) 2-Chlorophenol	6.67	128	468415	71.59	ng	98
10) Phenol	6.54	94	858799	102.69	ng	68
17) 2-Methylphenol	7.13	107	424205	79.97	ng	# 85
20) 3+4-Methylphenols	7.29	107	967669	144.42	ng	86
26) 2-Nitrophenol	7.79	139	327004	95.57	ng	# 81
27) 2,4-Dimethylphenol	7.83	122	527304	87.60	ng	# 82
29) 2,4-Dichlorophenol	8.02	162	722045	123.71	ng	95
36) 4-Chloro-3-methylphenol	8.72	107	753915	132.10	ng	# 82
42) 2,4,6-Trichlorophenol	9.17	196	154603	37.53	ng	97
43) 2,4,5-Trichlorophenol	9.20	196	179260m	42.32	ng	
53) 2,4-Dinitrophenol	9.99	184	191025	107.58	ng	# 48
55) 4-Nitrophenol	10.05	139	304777	124.30	ng	# 51
64) 4,6-Dinitro-2-methylphenol	10.53	198	114886	59.81	ng	# 59
69) Pentachlorophenol	11.22	266	344672	132.81	ng	99

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

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