

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100515\
 Data File : BF082049.D
 Acq On : 5 Oct 2015 16:54
 Operator : UM/IZ
 Sample : G3692-06DL 5X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PT-ACIDS-WPDL

Manual Integrations
 APPROVED

MMDadoda
 10/6/2015 2:55:23 PM

Quant Time: Oct 05 17:30:31 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	90906m	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	391160	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	178475	20.00	ng	-0.01
63) Phenanthrene-d10	11.42	188	355390	20.00	ng	0.00
75) Chrysene-d12	14.06	240	338011	20.00	ng	0.00
86) Perylene-d12	15.50	264	297673	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	143999	28.76	ng	-0.01
7) Phenol-d6	6.50	99	201728m	32.01	ng	-0.01
23) Nitrobenzene-d5	7.45	82	123189	20.99	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	59079	32.69	ng	-0.01
44) 2-Fluorobiphenyl	9.25	172	292314	23.46	ng	0.00
78) Terphenyl-d14	13.00	244	299472	22.73	ng	-0.01
Target Compounds						Qvalue
8) 2-Chlorophenol	6.66	128	97000	17.54	ng	100
10) Phenol	6.51	94	204367	28.91	ng	71
17) 2-Methylphenol	7.13	107	87730m	19.57	ng	
20) 3+4-Methylphenols	7.28	107	200036	35.32	ng	85
26) 2-Nitrophenol	7.78	139	58536	19.14	ng	# 51
27) 2,4-Dimethylphenol	7.82	122	102711	19.09	ng	# 78
29) 2,4-Dichlorophenol	8.02	162	163265	31.29	ng	98
36) 4-Chloro-3-methylphenol	8.71	107	167854	32.90	ng	# 74
42) 2,4,6-Trichlorophenol	9.15	196	26532	7.06	ng	97
43) 2,4,5-Trichlorophenol	9.20	196	32573	8.43	ng	97
53) 2,4-Dinitrophenol	9.99	184	25927	33.13	ng	# 67
55) 4-Nitrophenol	10.05	139	59167	26.46	ng	# 66
64) 4,6-Dinitro-2-methylphenol	10.51	198	17042	13.34	ng	89
69) Pentachlorophenol	11.21	266	62261	27.02	ng	97

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

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