

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080315\
 Data File : BF080818.D
 Acq On : 4 Aug 2015 00:06
 Operator : TP
 Sample : G3083-12
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 8535

Manual Integrations
 APPROVED

MMDadoda
 8/5/2015 2:28:23 PM

Quant Time: Aug 04 04:07:43 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 04 01:58:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	166376	20.00	ng	-0.01
21) Naphthalene-d8	8.11	136	582586	20.00	ng	-0.01
38) Acenaphthene-d10	9.87	164	283043	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	474309	20.00	ng	0.00
75) Chrysene-d12	13.97	240	323991	20.00	ng	-0.01
86) Perylene-d12	15.39	264	256931	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	937765	93.97	ng	0.00
7) Phenol-d6	6.46	99	853479	63.58	ng	0.00
23) Nitrobenzene-d5	7.40	82	1286513	108.97	ng	0.00
41) 2,4,6-Tribromophenol	10.66	330	475340	175.26	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	1970646	105.24	ng	0.00
78) Terphenyl-d14	12.93	244	1431329	88.04	ng	0.00
Target Compounds						Qvalue
8) 2-Chlorophenol	6.61	128	526350	47.61	ng	# 68
10) Phenol	6.48	94	842395	53.73	ng	94
17) 2-Methylphenol	7.08	107	334954	34.36	ng	98
20) 3+4-Methylphenols	7.24	107	639278	56.46	ng	# 57
26) 2-Nitrophenol	7.74	139	878102	170.48	ng	# 70
27) 2,4-Dimethylphenol	7.77	122	488949	49.82	ng	95
29) 2,4-Dichlorophenol	7.97	162	275169	33.51	ng	90
36) 4-Chloro-3-methylphenol	8.67	107	1558565	178.21	ng	89
42) 2,4,6-Trichlorophenol	9.10	196	793410	124.31	ng	98
43) 2,4,5-Trichlorophenol	9.14	196	467788m	74.89	ng	
53) 2,4-Dinitrophenol	9.93	184	316416	151.86	ng	98
55) 4-Nitrophenol	9.97	139	239556	70.91	ng	94
69) Pentachlorophenol	11.15	266	301847	95.85	ng	97

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

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