

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101415\
 Data File : BM002903.D
 Acq On : 14 Oct 2015 18:52
 Operator : UM/IZ
 Sample : G3692-08 25X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PT-ACIDS-WP

Quant Time: Oct 15 01:04:01 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-ACID-BM0101315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 15 01:02:42 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.66	152	83489	5.00	ng	0.01
8) Naphthalene-d8	11.50	136	339719	5.00	ng	0.00
14) Acenaphthene-d10	15.25	164	173610	5.00	ng	0.00
21) Phenanthrene-d10	17.95	188	408417	5.00	ng	0.00
24) Chrysene-d12	22.04	240	382314	5.00	ng	0.00
26) Perylene-d12	24.61	264	351167	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	6.12	112	93355	5.80	ng	-0.03
3) Phenol-d6	7.77	99	122862	5.36	ng	-0.03
9) Nitrobenzene-d5	9.85	82	66337	3.54	ng	-0.01
15) 2,4,6-Tribromophenol	16.71	330	39436	6.22	ng	0.00
18) 2-Fluorobiphenyl	13.87	172	224744	3.96	ng	0.00
25) Terphenyl-d14	20.48	244	217023	3.73	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) 2-Chlorophenol	8.20	128	64023	3.01	ng	# 71
5) Phenol	7.80	94	121111	5.38	ng	90
6) 2-Methylphenol	9.10	107	54936	3.19	ng	# 89
7) 3+4-Methylphenols	9.41	107	131881	6.19	ng	# 29
10) 2-Nitrophenol	10.62	139	28996	3.39	ng	# 74
11) 2,4-Dimethylphenol	10.64	122	68683	3.66	ng	91
12) 2,4-Dichlorophenol	11.17	162	88568	6.06	ng	97
13) 4-Chloro-3-methylphenol	12.74	107	119372	6.53	ng	87
16) 2,4,6-Trichlorophenol	13.72	196	16636	1.47	ng	# 73
17) 2,4,5-Trichlorophenol	13.83	196	19495	1.62	ng	# 92
19) 2,4-Dinitrophenol	15.43	184	4897	3.89	ng	96
20) 4-Nitrophenol	15.47	139	24017	3.38	ng	# 70
22) 4,6-Dinitro-2-methylphenol	16.38	198	5699	1.81	ng	# 73
23) Pentachlorophenol	17.62	266	18852	4.18	ng	91

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

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