

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060415\
 Data File : BG017468.D
 Acq On : 5 Jun 2015 8:52
 Operator : TP/IZ
 Sample : G2474-01
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C-1

Manual Integrations
 APPROVED

apatel
 6/5/2015 5:15:38 PM

Quant Time: Jun 05 15:27:19 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 04 01:31:15 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	21288	20.00	ng	0.00
21) Naphthalene-d8	10.42	136	91157	20.00	ng	0.00
38) Acenaphthene-d10	14.29	164	63928	20.00	ng	0.00
63) Phenanthrene-d10	17.04	188	149742	20.00	ng	0.00
75) Chrysene-d12	21.24	240	165004	20.00	ng	0.00
86) Perylene-d12	23.47	264	165315	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	140867	116.58	ng	0.01
7) Phenol-d6	6.85	99	197204	119.80	ng	0.01
23) Nitrobenzene-d5	8.79	82	120755	66.13	ng	0.00
41) 2,4,6-Tribromophenol	15.79	330	84528	95.20	ng	0.00
44) 2-Fluorobiphenyl	12.90	172	239328	56.47	ng	0.00
78) Terphenyl-d14	19.68	244	351479	44.03	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	13.75	163	33502	6.02	ng	98
70) Phenanthrene	17.08	178	32782	3.93	ng	96
74) Fluoranthene	19.11	202	68750	6.30	ng	98
77) Pyrene	19.47	202	65657	6.46	ng	99
80) Benzo(a)anthracene	21.22	228	35980	3.55	ng	97
82) Chrysene	21.27	228	38170	4.15	ng	95
83) Bis(2-ethylhexyl)phthalate	21.15	149	20325	3.20	ng	98
85) Indeno(1,2,3-cd)pyrene	25.75	276	26605	2.31	ng	# 79
87) Benzo(b)fluoranthene	22.79	252	46285	4.42	ng	# 89
88) Benzo(k)fluoranthene	22.83	252	20541m	2.05	ng	
89) Benzo(a)pyrene	23.37	252	33793	3.52	ng	93
91) Benzo(a,h,i)perylene	26.44	276	24014	2.53	ng	99

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

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