

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061316\
 Data File : BG022330.D
 Acq On : 13 Jun 2016 18:57
 Operator : UM/SJ
 Sample : H3474-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S7-B4(6-12)C

Quant Time: Jun 14 10:50:36 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 13 17:40:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.09	152	37668	20.00	ng	0.00
21) Naphthalene-d8	10.91	136	171667	20.00	ng	0.00
38) Acenaphthene-d10	14.72	164	91204	20.00	ng	0.00
63) Phenanthrene-d10	17.47	188	179700	20.00	ng	0.00
75) Chrysene-d12	21.74	240	147269	20.00	ng	0.00
86) Perylene-d12	25.02	264	130612	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	261177	134.06	ng	0.00
7) Phenol-d6	7.25	99	372695	121.67	ng	0.00
23) Nitrobenzene-d5	9.26	82	203668	82.71	ng	0.00
41) 2,4,6-Tribromophenol	16.21	330	106919	103.75	ng	0.00
44) 2-Fluorobiphenyl	13.35	172	430036	71.86	ng	0.00
78) Terphenyl-d14	20.06	244	413891	66.72	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.17	163	76658	10.14	ng	99
70) Phenanthrene	17.51	178	78888	8.08	ng	98
71) Anthracene	17.60	178	24676	2.55	ng	93
74) Fluoranthene	19.52	202	353574	31.51	ng	96
77) Pyrene	19.88	202	289967	31.67	ng	99
80) Benzo(a)anthracene	21.72	228	192261	23.28	ng	97
82) Chrysene	21.79	228	163618	20.36	ng	97
85) Indeno(1,2,3-cd)pyrene	28.80	276	52368	5.81	ng	# 54
87) Benzo(b)fluoranthene	23.99	252	178789m	23.47	ng	
88) Benzo(k)fluoranthene	24.05	252	96806m	12.91	ng	
89) Benzo(a)pyrene	24.87	252	128015	17.58	ng	# 93
91) Benzo(a,h,i)perylene	29.96	276	43882m	6.15	ng	

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

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