

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022916\
 Data File : BG021182.D
 Acq On : 1 Mar 2016 10:46
 Operator : UM/SJ
 Sample : H1520-19RE
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 40050+15(0-2)RE

Manual Integrations
 APPROVED

Sohil
 3/1/2016 6:51:11 PM

Quant Time: Mar 01 14:29:16 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 29 17:40:25 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	13803	20.00	ng	0.00
21) Naphthalene-d8	10.96	136	66153	20.00	ng	0.00
38) Acenaphthene-d10	14.76	164	36384	20.00	ng	0.00
63) Phenanthrene-d10	17.49	188	76211	20.00	ng	0.00
75) Chrysene-d12	21.76	240	80620	20.00	ng	0.00
86) Perylene-d12	25.03	264	74692	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.71	112	123945	142.75	ng	0.00
7) Phenol-d6	7.30	99	188401	131.96	ng	0.00
23) Nitrobenzene-d5	9.31	82	135395	86.61	ng	0.00
41) 2,4,6-Tribromophenol	16.24	330	45719	120.77	ng	0.00
44) 2-Fluorobiphenyl	13.38	172	207289	80.90	ng	0.00
78) Terphenyl-d14	20.08	244	245790	73.86	ng	0.00
Target Compounds						Qvalue
20) 3+4-Methylphenols	8.90	107	4057	2.72	ng	91
49) Dimethylphthalate	14.20	163	44609	13.76	ng	96
51) Acenaphthene	14.82	154	5786	2.40	ng	94
70) Phenanthrene	17.53	178	72282	17.57	ng	98
71) Anthracene	17.63	178	17718	4.21	ng	99
74) Fluoranthene	19.53	202	102450	21.20	ng	99
77) Pyrene	19.89	202	81359	16.92	ng	99
80) Benzo(a)anthracene	21.73	228	41454	8.68	ng	98
82) Chrysene	21.80	228	34573	7.64	ng	96
85) Indeno(1,2,3-cd)pyrene	28.75	276	15135	2.92	ng	# 100
87) Benzo(b)fluoranthene	23.99	252	44955	9.57	ng	# 97
88) Benzo(k)fluoranthene	24.05	252	18212m	3.87	ng	
89) Benzo(a)pyrene	24.87	252	31021	6.83	ng	# 88
91) Benzo(a,h,i)perylene	29.92	276	13112	3.02	ng	93

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

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