

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
 Data File : BG024056.D
 Acq On : 21 Sep 2016 15:48
 Operator : UM/SJ
 Sample : H4819-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YORK-SB2-02

Quant Time: Sep 21 16:34:28 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG083116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 19:01:00 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	79650	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	324846	20.00	ng	0.00
38) Acenaphthene-d10	14.72	164	231662	20.00	ng	0.00
63) Phenanthrene-d10	17.49	188	433389	20.00	ng	0.00
75) Chrysene-d12	21.80	240	458850	20.00	ng	0.00
86) Perylene-d12	25.15	264	445809	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	357273	78.75	ng	0.00
7) Phenol-d6	7.22	99	558639	79.99	ng	0.00
23) Nitrobenzene-d5	9.23	82	431378	59.28	ng	0.00
41) 2,4,6-Tribromophenol	16.23	330	216755	51.93	ng	0.00
44) 2-Fluorobiphenyl	13.33	172	803879	49.75	ng	0.00
78) Terphenyl-d14	20.10	244	1000275	49.65	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	10.93	128	76317	4.56	ng	# 93
49) Dimethylphthalate	14.17	163	191308	9.63	ng	99
70) Phenanthrene	17.53	178	332062	14.73	ng	96
71) Anthracene	17.62	178	63746	2.77	ng	95
74) Fluoranthene	19.55	202	636739	23.46	ng	94
77) Pyrene	19.91	202	524905	18.60	ng	94
80) Benzo(a)anthracene	21.78	228	358619	13.96	ng	97
82) Chrysene	21.85	228	372021	15.22	ng	99
85) Indeno(1,2,3-cd)pyrene	28.99	276	132047	4.88	ng	# 100
87) Benzo(b)fluoranthene	24.08	252	444487	17.55	ng	# 84
88) Benzo(k)fluoranthene	24.14	252	143918m	5.73	ng	
89) Benzo(a)pyrene	24.99	252	265714	11.05	ng	# 94
91) Benzo(a,h,i)perylene	30.18	276	46616	2.05	ng	# 89

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

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