

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
 Data File : BG024061.D
 Acq On : 21 Sep 2016 18:58
 Operator : UM/SJ
 Sample : H4820-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YORK-TP1-03

Quant Time: Sep 22 07:34:10 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	84730	20.00	ng	0.00
21) Naphthalene-d8	10.89	136	344082	20.00	ng	0.00
38) Acenaphthene-d10	14.73	164	237425	20.00	ng	0.00
63) Phenanthrene-d10	17.49	188	424580	20.00	ng	0.00
75) Chrysene-d12	21.80	240	437022	20.00	ng	0.00
86) Perylene-d12	25.15	264	422756	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	468789	97.14	ng	0.00
7) Phenol-d6	7.23	99	713322	96.02	ng	0.00
23) Nitrobenzene-d5	9.24	82	544683	70.67	ng	0.00
41) 2,4,6-Tribromophenol	16.23	330	270055	63.13	ng	0.00
44) 2-Fluorobiphenyl	13.34	172	952282	57.50	ng	0.00
78) Terphenyl-d14	20.10	244	1146731	59.76	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.17	163	259583	12.75	ng	97
70) Phenanthrene	17.54	178	104246	4.72	ng	95
74) Fluoranthene	19.55	202	212962	8.01	ng	91
77) Pyrene	19.92	202	234677	8.73	ng	93
80) Benzo(a)anthracene	21.78	228	158114m	6.46	ng	
82) Chrysene	21.85	228	148228	6.37	ng	98
85) Indeno(1,2,3-cd)pyrene	28.99	276	104474	4.05	ng	# 100
87) Benzo(b)fluoranthene	24.09	252	207203	8.63	ng	# 98
88) Benzo(k)fluoranthene	24.14	252	80381m	3.38	ng	
89) Benzo(a)pyrene	24.99	252	159471	6.99	ng	# 96
91) Benzo(g,h,i)perylene	30.21	276	103091	4.78	ng	# 92

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

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