

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
 Data File : BG022239.D
 Acq On : 8 Jun 2016 17:30
 Operator : UM/SJ
 Sample : H3402-02 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SS-39

Quant Time: Jun 09 06:45:13 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 06 16:12:45 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	39701	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	203813	20.00	ng	0.00
38) Acenaphthene-d10	14.75	164	121273	20.00	ng	0.00
63) Phenanthrene-d10	17.49	188	236572	20.00	ng	0.00
75) Chrysene-d12	21.76	240	210990	20.00	ng	0.00
86) Perylene-d12	25.05	264	205512	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	98732	48.08	ng	0.00
7) Phenol-d6	7.27	99	151376	46.89	ng	0.00
23) Nitrobenzene-d5	9.29	82	78449	26.83	ng	0.00
41) 2,4,6-Tribromophenol	16.23	330	49268	35.95	ng	0.00
44) 2-Fluorobiphenyl	13.37	172	232430	29.21	ng	0.00
78) Terphenyl-d14	20.08	244	253466	28.52	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.19	163	31887	3.17	ng	# 96
70) Phenanthrene	17.53	178	130226	10.14	ng	97
71) Anthracene	17.63	178	34337	2.69	ng	100
74) Fluoranthene	19.54	202	224394	15.19	ng	96
77) Pyrene	19.89	202	205383	15.66	ng	99
80) Benzo(a)anthracene	21.74	228	105843	8.95	ng	97
82) Chrysene	21.81	228	83247	7.23	ng	96
85) Indeno(1,2,3-cd)pyrene	28.83	276	54226	4.20	ng	# 88
87) Benzo(b)fluoranthene	24.01	252	126985m	10.59	ng	
88) Benzo(k)fluoranthene	24.07	252	31863m	2.70	ng	
89) Benzo(a)pyrene	24.91	252	84297	7.36	ng	97
91) Benzo(a,h,i)perylene	30.02	276	55255	4.92	ng	97

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

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