

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022316\
 Data File : BG021044.D
 Acq On : 24 Feb 2016 3:15
 Operator : UM/SJ
 Sample : H1507-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P-01

Manual Integrations
 APPROVED

Sohil
 2/24/2016 4:29:39 PM

Quant Time: Feb 24 04:39:42 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 24 00:35:06 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.21	152	4937	20.00	ng	0.00
21) Naphthalene-d8	11.03	136	26541	20.00	ng	0.00
38) Acenaphthene-d10	14.82	164	15402	20.00	ng	0.00
63) Phenanthrene-d10	17.55	188	27120	20.00	ng	0.00
75) Chrysene-d12	21.82	240	17532	20.00	ng	0.00
86) Perylene-d12	25.14	264	15962	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.79	112	46130	163.81	ng	0.00
7) Phenol-d6	7.41	99	74204	171.10	ng	0.00
23) Nitrobenzene-d5	9.39	82	48444	107.17	ng	0.00
41) 2,4,6-Tribromophenol	16.31	330	28637	155.20	ng	0.00
44) 2-Fluorobiphenyl	13.45	172	120530	108.94	ng	0.00
78) Terphenyl-d14	20.14	244	95350	70.57	ng	0.00
Target Compounds						
49) Dimethylphthalate	14.28	163	20281	16.15	ng	Qvalue # 98
70) Phenanthrene	17.59	178	34260	22.33	ng	98
71) Anthracene	17.69	178	10831	7.09	ng	98
74) Fluoranthene	19.59	202	53544	34.71	ng	99
77) Pyrene	19.95	202	43148	21.54	ng	99
80) Benzo(a)anthracene	21.80	228	17257	16.42	ng	98
82) Chrysene	21.87	228	14830	15.12	ng	98
87) Benzo(b)fluoranthene	24.08	252	14326	13.61	ng	97
88) Benzo(k)fluoranthene	24.14	252	5236m	5.27	ng	
89) Benzo(a)pyrene	24.98	252	8620	9.14	ng	# 93
91) Benzo(g,h,i)perylene	30.09	276	4723	5.57	ng	95

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

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