

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072115\
 Data File : BG018039.D
 Acq On : 22 Jul 2015 2:32
 Operator : TP/UM
 Sample : G3031-05
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-2(3-5)

Quant Time: Jul 22 05:16:17 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 22 04:14:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	61475	20.00	ng	0.00
21) Naphthalene-d8	10.33	136	281017	20.00	ng	0.00
38) Acenaphthene-d10	14.21	164	174739	20.00	ng	0.00
63) Phenanthrene-d10	16.96	188	359671	20.00	ng	0.00
75) Chrysene-d12	21.16	240	382534	20.00	ng	0.00
86) Perylene-d12	23.36	264	386439	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	491354	139.43	ng	0.00
7) Phenol-d6	6.74	99	711290	154.06	ng	0.00
23) Nitrobenzene-d5	8.71	82	424940	97.94	ng	0.00
41) 2,4,6-Tribromophenol	15.71	330	342311	193.36	ng	0.00
44) 2-Fluorobiphenyl	12.82	172	1156903	114.73	ng	0.00
78) Terphenyl-d14	19.62	244	1543008	101.61	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	13.67	163	175440	14.45	ng	98
70) Phenanthrene	17.00	178	67464	3.64	ng	98
74) Fluoranthene	19.03	202	229057	11.28	ng	95
77) Pyrene	19.39	202	225685	10.58	ng	95
80) Benzo(a)anthracene	21.15	228	140317	6.95	ng	98
82) Chrysene	21.20	228	155664m	8.09	ng	
83) Bis(2-ethylhexyl)phthalate	21.10	149	121774	9.26	ng	98
85) Indeno(1,2,3-cd)pyrene	25.59	276	135827m	5.84	ng	
87) Benzo(b)fluoranthene	22.71	252	254479m	12.04	ng	
88) Benzo(k)fluoranthene	22.74	252	66288m	3.14	ng	
89) Benzo(a)pyrene	23.27	252	165623	8.41	ng	99
91) Benzo(g,h,i)perylene	26.27	276	125107	6.35	ng	97

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

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