

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072115\
 Data File : BG018026.D
 Acq On : 21 Jul 2015 17:09
 Operator : TP/UM
 Sample : G2954-20
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TEC-B2

Quant Time: Jul 22 11:31:44 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 01:57:03 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	137407	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	623085	20.00	ng	0.00
38) Acenaphthene-d10	14.23	164	379583	20.00	ng	0.00
63) Phenanthrene-d10	16.99	188	726392	20.00	ng	0.00
75) Chrysene-d12	21.20	240	792957	20.00	ng	0.01
86) Perylene-d12	23.42	264	834092	20.00	ng	0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.18	112	919442	116.73	ng	0.00
7) Phenol-d6	6.76	99	1237756	119.94	ng	0.01
23) Nitrobenzene-d5	8.73	82	734217	76.32	ng	0.00
41) 2,4,6-Tribromophenol	15.73	330	469348	122.04	ng	0.00
44) 2-Fluorobiphenyl	12.85	172	1384884	63.22	ng	0.00
78) Terphenyl-d14	19.64	244	1588022	50.45	ng	0.00

Target Compounds						Qvalue
31) Naphthalene	10.40	128	100627	3.30	ng	98
48) Acenaphthylene	13.95	152	121421	3.38	ng	97
49) Dimethylphthalate	13.69	163	308730	11.71	ng	99
51) Acenaphthene	14.29	154	101315	4.66	ng	97
54) Dibenzofuran	14.63	168	142271	4.53	ng	# 95
57) Fluorene	15.28	166	183236	6.79	ng	98
70) Phenanthrene	17.03	178	2336992m	62.37	ng	
71) Anthracene	17.12	178	818553	22.45	ng	98
72) Carbazole	17.39	167	118243	3.43	ng	98
74) Fluoranthene	19.07	202	3154029	76.89	ng	81
77) Pyrene	19.43	202	3114874	70.43	ng	# 78
80) Benzo(a)anthracene	21.18	228	1890393	45.14	ng	# 90
82) Chrysene	21.24	228	1682999m	42.19	ng	
85) Indeno(1,2,3-cd)pyrene	25.68	276	936046	19.40	ng	# 85
87) Benzo(b)fluoranthene	22.76	252	2243264m	49.15	ng	
88) Benzo(k)fluoranthene	22.79	252	980861m	21.52	ng	
89) Benzo(a)pyrene	23.33	252	1920781	45.18	ng	95
90) Dibenzo(a,h)anthracene	25.68	278	224232m	5.09	ng	
91) Benzo(g,h,i)perylene	26.38	276	967658	22.76	ng	94

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

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