

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042815\
 Data File : BG016779.D
 Acq On : 28 Apr 2015 23:40
 Operator : TP/IZ
 Sample : G2003-01 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-18(40-42)

Quant Time: Apr 29 02:01:42 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 29 00:55:31 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	72778	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	304050	20.00	ng	0.00
38) Acenaphthene-d10	14.22	164	159520	20.00	ng	0.00
63) Phenanthrene-d10	16.97	188	293800	20.00	ng	0.00
75) Chrysene-d12	21.16	240	251291	20.00	ng	0.00
86) Perylene-d12	23.36	264	247536	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	186475	41.58	ng	0.00
7) Phenol-d6	6.77	99	256351	40.74	ng	0.00
23) Nitrobenzene-d5	8.73	82	131132	27.25	ng	0.00
41) 2,4,6-Tribromophenol	15.72	330	48563	41.04	ng	0.00
44) 2-Fluorobiphenyl	12.84	172	300565	30.27	ng	0.00
78) Terphenyl-d14	19.60	244	291053	26.68	ng	0.00

Target Compounds				Qvalue		
48) Acenaphthylene	13.95	152	52786	3.13	ng	99
54) Dibenzofuran	14.62	168	48023	3.36	ng	# 97
57) Fluorene	15.27	166	76248	6.11	ng	96
70) Phenanthrene	17.01	178	772825	45.79	ng	98
71) Anthracene	17.10	178	131407	7.83	ng	99
72) Carbazole	17.38	167	40293	2.40	ng	95
74) Fluoranthene	19.03	202	786533	43.98	ng	98
77) Pyrene	19.40	202	613041	34.73	ng	100
80) Benzo(a)anthracene	21.14	228	310818	20.40	ng	99
82) Chrysene	21.20	228	265693	18.19	ng	98
85) Indeno(1,2,3-cd)pyrene	25.58	276	98260	6.17	ng	91
87) Benzo(b)fluoranthene	22.69	252	291418	19.38	ng	99
88) Benzo(k)fluoranthene	22.74	252	108892m	7.42	ng	
89) Benzo(a)pyrene	23.26	252	197818	13.96	ng	97
90) Dibenzo(a,h)anthracene	25.58	278	28787	2.09	ng	96
91) Benzo(g,h,i)perylene	26.27	276	79158	5.65	ng	96

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

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