

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042815\
 Data File : BG016778.D
 Acq On : 28 Apr 2015 23:05
 Operator : TP/IZ
 Sample : G2022-04 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-04-COMP

Quant Time: Apr 29 01:59:12 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	71731	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	300209	20.00	ng	0.00
38) Acenaphthene-d10	14.22	164	162692	20.00	ng	0.00
63) Phenanthrene-d10	16.97	188	298172	20.00	ng	0.00
75) Chrysene-d12	21.16	240	247470	20.00	ng	0.00
86) Perylene-d12	23.36	264	242605	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	223779	50.62	ng	0.00
7) Phenol-d6	6.76	99	304795	49.15	ng	0.00
23) Nitrobenzene-d5	8.73	82	158563	33.37	ng	0.00
41) 2,4,6-Tribromophenol	15.72	330	59954	49.67	ng	0.00
44) 2-Fluorobiphenyl	12.84	172	352899	34.85	ng	0.00
78) Terphenyl-d14	19.60	244	326086	30.35	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	10.40	128	166699	10.53	ng	96
37) 2-Methylnaphthalene	12.02	142	95676	8.37	ng	100
70) Phenanthrene	17.01	178	145009	8.47	ng	98
74) Fluoranthene	19.03	202	245005	13.50	ng	97
77) Pyrene	19.40	202	247102	14.21	ng	97
80) Benzo(a)anthracene	21.14	228	124856	8.32	ng	99
82) Chrysene	21.19	228	116686	8.11	ng	96
83) Bis(2-ethylhexyl)phthalate	21.08	149	155862	13.40	ng	100
85) Indeno(1,2,3-cd)pyrene	25.58	276	67989	4.34	ng	# 85
87) Benzo(b)fluoranthene	22.70	252	149685	10.16	ng	96
88) Benzo(k)fluoranthene	22.73	252	55523m	3.86	ng	
89) Benzo(a)pyrene	23.26	252	116636	8.40	ng	94
91) Benzo(a,h,i)perylene	26.27	276	62877	4.58	ng	96

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

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