

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042915\
 Data File : BF078834.D
 Acq On : 30 Apr 2015 3:15
 Operator : TP/IZ
 Sample : G2022-14
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-10-COMP

Quant Time: Apr 30 04:07:13 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 30 01:56:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	67935	20.00	ng	0.00
21) Naphthalene-d8	8.96	136	289214	20.00	ng	0.00
38) Acenaphthene-d10	11.14	164	136556	20.00	ng	0.00
63) Phenanthrene-d10	12.98	188	292424	20.00	ng	0.00
75) Chrysene-d12	16.27	240	323299	20.00	ng	-0.01
86) Perylene-d12	17.99	264	357468	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.68	112	318336	86.06	ng	0.01
7) Phenol-d6	6.92	99	457445	92.60	ng	0.00
23) Nitrobenzene-d5	8.07	82	240728	59.93	ng	0.00
41) 2,4,6-Tribromophenol	12.11	330	120792	90.63	ng	-0.01
44) 2-Fluorobiphenyl	10.31	172	511900	51.34	ng	0.00
78) Terphenyl-d14	14.98	244	653838	48.16	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.94	94	18592	3.18	ng	94
49) Dimethylphthalate	10.81	163	27061	2.79	ng	99
70) Phenanthrene	13.02	178	46112	2.91	ng	97
74) Fluoranthene	14.49	202	98607	6.25	ng	98
77) Pyrene	14.78	202	95750	5.07	ng	99
80) Benzo(a)anthracene	16.26	228	61508	3.55	ng	94
82) Chrysene	16.31	228	48568	2.87	ng	96
85) Indeno(1,2,3-cd)pyrene	19.50	276	47730	2.49	ng	# 100
87) Benzo(b)fluoranthene	17.54	252	76266m	3.94	ng	
89) Benzo(a)pyrene	17.92	252	57352	3.31	ng	# 97
91) Benzo(g,h,i)perylene	19.94	276	51218	2.85	ng	96

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

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