

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
 Data File : BF079171.D
 Acq On : 12 May 2015 21:40
 Operator : TP/IZ
 Sample : G2175-14
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-H

Quant Time: May 13 02:17:22 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 13 00:53:14 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.28	152	49664	20.00	ng	0.00
21) Naphthalene-d8	8.86	136	209858	20.00	ng	0.00
38) Acenaphthene-d10	11.03	164	100907	20.00	ng	-0.01
63) Phenanthrene-d10	12.88	188	194438	20.00	ng	0.00
75) Chrysene-d12	16.16	240	213237	20.00	ng	-0.03
86) Perylene-d12	17.86	264	227780	20.00	ng	-0.16
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	113394	39.30	ng	0.01
7) Phenol-d6	6.84	99	179193	46.99	ng	0.00
23) Nitrobenzene-d5	7.98	82	93957	25.73	ng	0.00
41) 2,4,6-Tribromophenol	12.01	330	117693	107.81	ng	0.00
44) 2-Fluorobiphenyl	10.20	172	315287	43.53	ng	0.00
78) Terphenyl-d14	14.87	244	585913	65.78	ng	-0.01
Target Compounds						Qvalue
32) Benzoic acid	8.51	122	2283	7.37	ng	98
48) Acenaphthylene	10.86	152	21484	2.09	ng	98
49) Dimethylphthalate	10.71	163	18801	2.54	ng	98
70) Phenanthrene	12.90	178	40971	3.77	ng	99
74) Fluoranthene	14.39	202	96936	8.55	ng	92
77) Pyrene	14.66	202	109834	8.94	ng	98
80) Benzo(a)anthracene	16.15	228	54081m	4.63	ng	
82) Chrysene	16.19	228	60331	5.37	ng	98
85) Indeno(1,2,3-cd)pyrene	19.31	276	38018m	2.66	ng	
87) Benzo(b)fluoranthene	17.43	252	74575	5.98	ng	98
88) Benzo(k)fluoranthene	17.44	252	32880m	2.72	ng	
89) Benzo(a)pyrene	17.79	252	54793	4.77	ng	# 91
91) Benzo(a,h,i)perylene	19.74	276	40054	3.28	ng	# 90

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

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