

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
 Data File : BF079170.D
 Acq On : 12 May 2015 21:11
 Operator : TP/IZ
 Sample : G2155-08 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-43AS

Quant Time: May 13 02:15:25 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	49960	20.00	ng	0.00
21) Naphthalene-d8	8.86	136	208521	20.00	ng	0.00
38) Acenaphthene-d10	11.03	164	102613	20.00	ng	-0.01
63) Phenanthrene-d10	12.88	188	203390	20.00	ng	0.00
75) Chrysene-d12	16.17	240	216251	20.00	ng	-0.02
86) Perylene-d12	17.90	264	223093	20.00	ng	-0.13
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	207891	71.63	ng	0.01
7) Phenol-d6	6.84	99	278139	72.50	ng	0.00
23) Nitrobenzene-d5	7.98	82	150087	41.37	ng	0.00
41) 2,4,6-Tribromophenol	12.01	330	76754	69.14	ng	0.00
44) 2-Fluorobiphenyl	10.20	172	320945	43.58	ng	0.00
78) Terphenyl-d14	14.87	244	376544	41.69	ng	-0.01
Target Compounds						Qvalue
32) Benzoic acid	8.51	122	1508	7.01	ng	93
49) Dimethylphthalate	10.71	163	22366	2.97	ng	99
70) Phenanthrene	12.90	178	66282	5.83	ng	97
74) Fluoranthene	14.39	202	97097	8.19	ng	91
77) Pyrene	14.66	202	106390	8.54	ng	97
80) Benzo(a)anthracene	16.16	228	52140	4.40	ng	96
82) Chrysene	16.19	228	50503	4.43	ng	97
83) Bis(2-ethylhexyl)phthalate	16.21	149	28196	3.42	ng	99
85) Indeno(1,2,3-cd)pyrene	19.35	276	35033m	2.41	ng	
87) Benzo(b)fluoranthene	17.45	252	69751	5.71	ng	96
89) Benzo(a)pyrene	17.83	252	48689	4.33	ng	# 88
91) Benzo(a,h,i)perylene	19.78	276	37260	3.11	ng	93

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

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