

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051815\
 Data File : BF079321.D
 Acq On : 18 May 2015 23:39
 Operator : TP/IZ
 Sample : G2270-03 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-19

Quant Time: May 19 01:58:29 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 19 01:26:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.16	152	44524	20.00	ng	0.00
21) Naphthalene-d8	8.74	136	183576	20.00	ng	0.00
38) Acenaphthene-d10	10.91	164	92372	20.00	ng	0.00
63) Phenanthrene-d10	12.75	188	172914	20.00	ng	0.00
75) Chrysene-d12	16.03	240	194857	20.00	ng	-0.02
86) Perylene-d12	17.77	264	203058	20.00	ng	-0.08
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	156487m	60.50	ng	0.00
7) Phenol-d6	6.73	99	204097	59.70	ng	-0.01
23) Nitrobenzene-d5	7.86	82	115662	36.21	ng	0.00
41) 2,4,6-Tribromophenol	11.90	330	60389	60.43	ng	0.00
44) 2-Fluorobiphenyl	10.09	172	247199	37.28	ng	0.00
78) Terphenyl-d14	14.75	244	295097	36.26	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	10.59	163	14351	2.12	ng	99
70) Phenanthrene	12.78	178	24375	2.52	ng	96
74) Fluoranthene	14.26	202	60571	6.01	ng	92
77) Pyrene	14.54	202	64886	5.78	ng	97
80) Benzo(a)anthracene	16.02	228	41271	3.87	ng	# 80
82) Chrysene	16.07	228	34279	3.34	ng	93
85) Indeno(1,2,3-cd)pyrene	19.15	276	30058	2.30	ng	96
87) Benzo(b)fluoranthene	17.33	252	51559m	4.64	ng	
88) Benzo(k)fluoranthene	17.34	252	22974m	2.14	ng	
89) Benzo(a)pyrene	17.70	252	42402	4.14	ng	# 96
91) Benzo(a,h,i)perylene	19.58	276	30001	2.75	ng	# 91

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

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