

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
 Data File : BF079159.D
 Acq On : 12 May 2015 15:55
 Operator : TP/IZ
 Sample : G2155-09
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-43D

Quant Time: May 13 01:30:37 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	69211	20.00	ng	0.00
21) Naphthalene-d8	8.86	136	292061	20.00	ng	0.00
38) Acenaphthene-d10	11.03	164	138895	20.00	ng	-0.01
63) Phenanthrene-d10	12.88	188	275586	20.00	ng	0.00
75) Chrysene-d12	16.21	240	276851	20.00	ng	0.01
86) Perylene-d12	18.09	264	298508	20.00	ng	0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	308592	76.75	ng	0.01
7) Phenol-d6	6.84	99	406864	76.56	ng	0.00
23) Nitrobenzene-d5	7.98	82	232198	45.69	ng	0.00
41) 2,4,6-Tribromophenol	12.01	330	115552	76.90	ng	0.00
44) 2-Fluorobiphenyl	10.20	172	440423	44.18	ng	0.00
78) Terphenyl-d14	14.87	244	457836	39.59	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	10.71	163	21328	2.09	ng	99
70) Phenanthrene	12.90	178	82410	5.35	ng	97
74) Fluoranthene	14.39	202	157565	9.81	ng	92
77) Pyrene	14.66	202	140001	8.78	ng	98
80) Benzo(a)anthracene	16.19	228	65119m	4.30	ng	
82) Chrysene	16.24	228	52933	3.63	ng	98
83) Bis(2-ethylhexyl)phthalate	16.26	149	67012	6.34	ng	100
85) Indeno(1,2,3-cd)pyrene	19.57	276	49045	2.64	ng	# 100
87) Benzo(b)fluoranthene	17.60	252	66426m	4.07	ng	
88) Benzo(k)fluoranthene	17.62	252	32017m	2.02	ng	
89) Benzo(a)pyrene	18.01	252	57220	3.80	ng	# 94
91) Benzo(a,h,i)perylene	20.00	276	43327	2.70	ng	98

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

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