

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051815\
 Data File : BF079328.D
 Acq On : 19 May 2015 3:00
 Operator : TP/IZ
 Sample : G2273-07
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-07-D1

Quant Time: May 19 04:55:06 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 04:38:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.16	152	45474	20.00	ng	0.00
21) Naphthalene-d8	8.74	136	189976	20.00	ng	0.00
38) Acenaphthene-d10	10.91	164	93576	20.00	ng	0.00
63) Phenanthrene-d10	12.75	188	184023	20.00	ng	0.00
75) Chrysene-d12	16.05	240	192985	20.00	ng	-0.01
86) Perylene-d12	17.76	264	207158	20.00	ng	-0.09

System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	280105m	106.03	ng	0.01
7) Phenol-d6	6.74	99	371893	106.50	ng	0.00
23) Nitrobenzene-d5	7.86	82	223847	67.72	ng	0.00
41) 2,4,6-Tribromophenol	11.90	330	113097	111.72	ng	0.00
44) 2-Fluorobiphenyl	10.09	172	451135	67.17	ng	0.00
78) Terphenyl-d14	14.75	244	526223	65.28	ng	0.00

Target Compounds						Qvalue
31) Naphthalene	8.76	128	20533	2.27	ng	99
48) Acenaphthylene	10.74	152	42553	4.47	ng	99
49) Dimethylphthalate	10.59	163	25448	3.71	ng	100
70) Phenanthrene	12.79	178	92840	9.02	ng	100
71) Anthracene	12.86	178	33473	3.31	ng	99
74) Fluoranthene	14.26	202	261609	24.39	ng	99
77) Pyrene	14.55	202	270555	24.33	ng	98
80) Benzo(a)anthracene	16.03	228	232487m	22.00	ng	
82) Chrysene	16.08	228	177418	17.45	ng	97
85) Indeno(1,2,3-cd)pyrene	19.15	276	170336	13.15	ng	96
87) Benzo(b)fluoranthene	17.33	252	305432m	26.94	ng	
88) Benzo(k)fluoranthene	17.35	252	124360m	11.33	ng	
89) Benzo(a)pyrene	17.69	252	237127	22.71	ng	# 95
90) Dibenzo(a,h)anthracene	19.18	278	42211	3.80	ng	# 83
91) Benzo(a,h,i)perylene	19.58	276	157364	14.15	ng	96

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

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