

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040615\
 Data File : BF078286.D
 Acq On : 7 Apr 2015 6:33
 Operator : TP/IZ
 Sample : G1779-01
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-2(1-3)

Quant Time: Apr 07 07:03:37 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 07 04:11:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	37533	20.00	ng	0.00
21) Naphthalene-d8	8.56	136	155085	20.00	ng	0.00
38) Acenaphthene-d10	10.72	164	78830	20.00	ng	0.00
63) Phenanthrene-d10	12.55	188	149726	20.00	ng	0.00
75) Chrysene-d12	15.81	240	165108	20.00	ng	0.00
86) Perylene-d12	17.51	264	163623	20.00	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	146477	64.35	ng	0.01
7) Phenol-d6	6.58	99	190201	66.67	ng	0.00
23) Nitrobenzene-d5	7.68	82	110186	43.06	ng	0.00
41) 2,4,6-Tribromophenol	11.70	330	42550	65.08	ng	0.00
44) 2-Fluorobiphenyl	9.90	172	233371	42.32	ng	0.00
78) Terphenyl-d14	14.55	244	271897	37.21	ng	0.00
Target Compounds						Qvalue
70) Phenanthrene	12.58	178	30959	3.57	ng	99
74) Fluoranthene	14.04	202	56726	6.21	ng	96
77) Pyrene	14.32	202	53385	5.15	ng	99
80) Benzo(a)anthracene	15.80	228	29010m	2.95	ng	
82) Chrysene	15.85	228	23894	2.59	ng	97
87) Benzo(b)fluoranthene	17.08	252	32170m	3.18	ng	
89) Benzo(a)pyrene	17.45	252	21804	2.47	ng	95

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

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