

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051915\
 Data File : BF079372.D
 Acq On : 20 May 2015 7:50
 Operator : TP/IZ
 Sample : G2289-06 5X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-33S

Quant Time: May 20 08:31:30 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 20 07:04:25 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.13	152	46194	20.00	ng	0.00
21) Naphthalene-d8	8.71	136	192722	20.00	ng	0.00
38) Acenaphthene-d10	10.87	164	91484	20.00	ng	-0.01
63) Phenanthrene-d10	12.72	188	177511	20.00	ng	0.00
75) Chrysene-d12	16.00	240	200330	20.00	ng	0.00
86) Perylene-d12	17.77	264	211318	20.00	ng	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.42	112	50664m	18.88	ng	0.01
7) Phenol-d6	6.70	99	64453m	18.17	ng	-0.01
23) Nitrobenzene-d5	7.82	82	35352	10.54	ng	-0.01
41) 2,4,6-Tribromophenol	11.85	330	16771	16.95	ng	-0.01
44) 2-Fluorobiphenyl	10.04	172	75363	11.48	ng	-0.01
78) Terphenyl-d14	14.71	244	87605	10.47	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
70) Phenanthrene	12.74	178	39656	3.99	ng	99
74) Fluoranthene	14.22	202	82646	7.99	ng	96
77) Pyrene	14.50	202	79423	6.88	ng	98
80) Benzo(a)anthracene	15.99	228	44530	4.06	ng	94
82) Chrysene	16.03	228	36548	3.46	ng	96
85) Indeno(1,2,3-cd)pyrene	19.15	276	27285	2.03	ng	# 90
87) Benzo(b)fluoranthene	17.31	252	48300m	4.18	ng	
89) Benzo(a)pyrene	17.70	252	38092	3.58	ng	# 91
91) Benzo(g,h,i)perylene	19.55	276	27296	2.41	ng	# 86

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

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