

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051915\
 Data File : BF079368.D
 Acq On : 20 May 2015 5:55
 Operator : TP/IZ
 Sample : G2289-02
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-34S

Quant Time: May 20 06:59:55 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 05:08:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.13	152	45619	20.00	ng	0.00
21) Naphthalene-d8	8.71	136	190088	20.00	ng	0.00
38) Acenaphthene-d10	10.88	164	93550	20.00	ng	0.00
63) Phenanthrene-d10	12.72	188	186123	20.00	ng	0.00
75) Chrysene-d12	16.00	240	200263	20.00	ng	-0.02
86) Perylene-d12	17.70	264	231135	20.00	ng	-0.11

System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	272081m	102.67	ng	0.00
7) Phenol-d6	6.71	99	356416	101.74	ng	0.00
23) Nitrobenzene-d5	7.82	82	200161	60.52	ng	-0.01
41) 2,4,6-Tribromophenol	11.86	330	106811	105.54	ng	0.00
44) 2-Fluorobiphenyl	10.06	172	426471	63.51	ng	0.00
78) Terphenyl-d14	14.71	244	484566	57.93	ng	-0.01

Target Compounds						Qvalue
31) Naphthalene	8.73	128	28201	3.11	ng	99
48) Acenaphthylene	10.70	152	23482	2.47	ng	100
49) Dimethylphthalate	10.56	163	26743	3.90	ng	100
51) Acenaphthene	10.91	154	49617	8.74	ng	99
54) Dibenzofuran	11.13	168	36569	4.46	ng	99
57) Fluorene	11.55	166	47436	7.04	ng	99
70) Phenanthrene	12.74	178	627989	60.31	ng	99
71) Anthracene	12.81	178	173697	17.00	ng	99
72) Carbazole	13.02	167	45298	4.65	ng	98
74) Fluoranthene	14.23	202	996218	91.82	ng	96
77) Pyrene	14.50	202	992557	86.01	ng	97
80) Benzo(a)anthracene	15.99	228	560389	51.10	ng	97
82) Chrysene	16.03	228	468628	44.43	ng	97
83) Bis(2-ethylhexyl)phthalate	16.05	149	19329	2.53	ng	# 91
85) Indeno(1,2,3-cd)pyrene	19.07	276	350326	26.07	ng	94
87) Benzo(b)fluoranthene	17.27	252	711113m	56.21	ng	
88) Benzo(k)fluoranthene	17.30	252	160129m	13.07	ng	
89) Benzo(a)pyrene	17.65	252	534755	45.91	ng	98
90) Dibenzo(a,h)anthracene	19.09	278	79946m	6.46	ng	
91) Benzo(a,h,i)perylene	19.49	276	355691	28.67	ng	97

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

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