

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030617\
 Data File : BF093405.D
 Acq On : 6 Mar 2017 17:18
 Operator : SJ/MA
 Sample : I2061-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 IMPACT-2-FROMDUMPSTER

Quant Time: Mar 07 00:13:06 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 07 00:03:56 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	175545	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	666314	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	222103	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	277476	20.00	ng	0.02
75) Chrysene-d12	13.93	240	358359	20.00	ng	0.00
86) Perylene-d12	15.36	264	325506	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	1595007	155.88	ng	0.02
7) Phenol-d6	6.42	99	1927057	146.37	ng	0.00
23) Nitrobenzene-d5	7.35	82	1274813	106.54	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	199913	124.45	ng	0.02
44) 2-Fluorobiphenyl	9.14	172	1646272	134.28	ng	0.00
78) Terphenyl-d14	12.88	244	988940	64.84	ng	0.00
Target Compounds						
10) Phenol	6.44	94	121438	7.88	ng	# 81
37) 2-Methylnaphthalene	8.78	142	50468	2.33	ng	95
49) Dimethylphthalate	9.53	163	236977	15.84	ng	# 94
77) Pyrene	12.75	202	67359	2.51	ng	# 91
83) Bis(2-ethylhexyl)phthalate	13.91	149	43310	2.91	ng	# 91
85) Indeno(1,2,3-cd)pyrene	16.76	276	38736	2.44	ng	98
91) Benzo(g,h,i)perylene	17.17	276	287644	19.01	ng	# 90

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

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