

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032217\
 Data File : BF093873.D
 Acq On : 23 Mar 2017 10:24
 Operator : SJ/MA
 Sample : I2337-05 2X
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-Q14-ESW

Quant Time: Mar 23 17:05:59 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 23 02:02:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.99	152	195282	20.00	ng	0.00
21) Naphthalene-d8	7.49	136	760278	20.00	ng	0.00
38) Acenaphthene-d10	9.63	164	324990	20.00	ng	0.00
63) Phenanthrene-d10	11.46	188	497400	20.00	ng	0.00
75) Chrysene-d12	14.71	240	340795	20.00	ng	0.00
86) Perylene-d12	16.34	264	319027	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.50	112	576766	49.88	ng	0.00
7) Phenol-d6	5.56	99	769394	53.79	ng	0.00
23) Nitrobenzene-d5	6.63	82	514514	38.62	ng	-0.01
41) 2,4,6-Tribromophenol	10.60	330	119329	46.47	ng	-0.01
44) 2-Fluorobiphenyl	8.81	172	770297	36.15	ng	-0.01
78) Terphenyl-d14	13.43	244	387946	25.52	ng	0.00
Target Compounds						Qvalue
10) Phenol	5.57	94	36591	2.25	ng	89
49) Dimethylphthalate	9.30	163	83960	3.63	ng	98
70) Phenanthrene	11.49	178	59242	2.14	ng	100
74) Fluoranthene	12.94	202	121925	4.30	ng	97
77) Pyrene	13.22	202	127513	5.16	ng	99
80) Benzo(a)anthracene	14.69	228	62131	3.13	ng	94
82) Chrysene	14.74	228	65053m	3.53	ng	
85) Indeno(1,2,3-cd)pyrene	17.73	276	57531	3.60	ng	97
87) Benzo(b)fluoranthene	15.93	252	104811m	5.36	ng	
89) Benzo(a)pyrene	16.28	252	82129	4.56	ng	98
91) Benzo(g,h,i)perylene	18.14	276	58363	3.80	ng	96

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

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