

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031517\
 Data File : BF093678.D
 Acq On : 15 Mar 2017 19:59
 Operator : SJ/MA
 Sample : I2264-09
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-DD8-BASE

Quant Time: Mar 16 03:04:38 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 13 15:05:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	166965	20.00	ng	-0.01
21) Naphthalene-d8	8.00	136	661825	20.00	ng	-0.01
38) Acenaphthene-d10	9.75	164	316484	20.00	ng	-0.01
63) Phenanthrene-d10	11.25	188	534642	20.00	ng	0.00
75) Chrysene-d12	13.89	240	320820	20.00	ng	0.00
86) Perylene-d12	15.29	264	304604	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	1194896	119.11	ng	-0.01
7) Phenol-d6	6.37	99	1547970	122.36	ng	-0.01
23) Nitrobenzene-d5	7.28	82	1012682	84.90	ng	-0.02
41) 2,4,6-Tribromophenol	10.55	330	243569	118.13	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	1349381	79.31	ng	-0.01
78) Terphenyl-d14	12.83	244	956783	77.54	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.38	94	126237	8.14	ng	87
49) Dimethylphthalate	9.49	163	400187	19.07	ng	100
70) Phenanthrene	11.27	178	74969	2.46	ng	97
74) Fluoranthene	12.47	202	145612	4.94	ng	92
77) Pyrene	12.69	202	136446	5.85	ng	98
80) Benzo(a)anthracene	13.88	228	69112	3.65	ng	99
82) Chrysene	13.92	228	51371	2.93	ng	98
85) Indeno(1,2,3-cd)pyrene	16.67	276	42281	2.88	ng	# 90
87) Benzo(b)fluoranthene	14.91	252	69730m	3.76	ng	
89) Benzo(a)pyrene	15.25	252	58407	3.44	ng	96
91) Benzo(g,h,i)perylene	17.08	276	47629	3.31	ng	# 89

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

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