

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031517\
 Data File : BF093677.D
 Acq On : 15 Mar 2017 19:31
 Operator : SJ/MA
 Sample : I2264-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-CC6-WSW

Quant Time: Mar 16 02:33:23 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	175817	20.00	ng	-0.01
21) Naphthalene-d8	8.00	136	710255	20.00	ng	-0.01
38) Acenaphthene-d10	9.76	164	329370	20.00	ng	0.00
63) Phenanthrene-d10	11.25	188	555542	20.00	ng	0.00
75) Chrysene-d12	13.89	240	359914	20.00	ng	0.00
86) Perylene-d12	15.28	264	319823	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	1167778	110.54	ng	-0.01
7) Phenol-d6	6.37	99	1545532	116.02	ng	-0.01
23) Nitrobenzene-d5	7.28	82	992571	77.54	ng	-0.02
41) 2,4,6-Tribromophenol	10.55	330	254274	118.49	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	1473920	83.24	ng	-0.01
78) Terphenyl-d14	12.83	244	1157184	83.59	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.38	94	118452	7.26	ng	91
49) Dimethylphthalate	9.49	163	301025	13.79	ng	99
70) Phenanthrene	11.27	178	136295	4.30	ng	96
74) Fluoranthene	12.46	202	203847	6.66	ng	97
77) Pyrene	12.69	202	187241	7.15	ng	98
80) Benzo(a)anthracene	13.88	228	86184	4.05	ng	99
82) Chrysene	13.91	228	67771	3.45	ng	98
85) Indeno(1,2,3-cd)pyrene	16.67	276	48419m	2.94	ng	
87) Benzo(b)fluoranthene	14.91	252	89688m	4.60	ng	
89) Benzo(a)pyrene	15.24	252	65859	3.70	ng	# 96
91) Benzo(g,h,i)perylene	17.07	276	49013	3.24	ng	# 84

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

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