

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031517\
 Data File : BF093683.D
 Acq On : 15 Mar 2017 22:16
 Operator : SJ/MA
 Sample : I2264-15
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-EE8-SSW

Quant Time: Mar 16 03:19:15 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	179336	20.00	ng	-0.01
21) Naphthalene-d8	8.00	136	650748	20.00	ng	-0.01
38) Acenaphthene-d10	9.75	164	255162	20.00	ng	-0.01
63) Phenanthrene-d10	11.25	188	367554	20.00	ng	0.00
75) Chrysene-d12	13.89	240	357458	20.00	ng	0.00
86) Perylene-d12	15.29	264	321907	20.00	ng	0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	1257093	116.66	ng	-0.01
7) Phenol-d6	6.37	99	1533249	112.84	ng	-0.01
23) Nitrobenzene-d5	7.28	82	934686	79.69	ng	-0.02
41) 2,4,6-Tribromophenol	10.55	330	170043	102.29	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	1201983	87.62	ng	-0.01
78) Terphenyl-d14	12.83	244	723138	52.60	ng	0.00

Target Compounds				Qvalue		
10) Phenol	6.39	94	119145	7.16	ng	80
31) Naphthalene	8.02	128	159665	4.90	ng	98
48) Acenaphthylene	9.62	152	53314	2.27	ng	98
49) Dimethylphthalate	9.49	163	151359	8.95	ng	100
51) Acenaphthene	9.78	154	53337	3.84	ng	97
54) Dibenzofuran	9.97	168	49376	2.84	ng	94
57) Fluorene	10.31	166	49902m	3.39	ng	
70) Phenanthrene	11.27	178	831593	39.64	ng	98
71) Anthracene	11.32	178	182708m	8.61	ng	
72) Carbazole	11.50	167	89545	4.49	ng	99
74) Fluoranthene	12.46	202	1093116	53.94	ng	97
77) Pvrene	12.69	202	1093865	42.08	ng	99
80) Benzo(a)anthracene	13.88	228	587038m	27.81	ng	
82) Chrysene	13.91	228	585994m	30.00	ng	
85) Indeno(1,2,3-cd)pyrene	16.65	276	284726	17.42	ng	98
87) Benzo(b)fluoranthene	14.91	252	725157m	36.99	ng	
88) Benzo(k)fluoranthene	14.93	252	232364m	12.29	ng	
89) Benzo(a)pyrene	15.24	252	533388	29.77	ng	# 92
90) Dibenzo(a,h)anthracene	16.64	278	64630	4.36	ng	# 89
91) Benzo(a,h,i)perylene	17.07	276	280976	18.47	ng	# 91

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

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