

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032217\
 Data File : BF093860.D
 Acq On : 23 Mar 2017 4:30
 Operator : SJ/MA
 Sample : I2337-07
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-Q14-BASE

Quant Time: Mar 23 07:02:17 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	196204	20.00	ng	0.00
21) Naphthalene-d8	7.49	136	788265	20.00	ng	0.00
38) Acenaphthene-d10	9.63	164	359923	20.00	ng	0.00
63) Phenanthrene-d10	11.46	188	619629	20.00	ng	0.00
75) Chrysene-d12	14.71	240	434685	20.00	ng	0.00
86) Perylene-d12	16.34	264	321559	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.52	112	1254491	107.99	ng	0.01
7) Phenol-d6	5.57	99	1608720	111.94	ng	0.00
23) Nitrobenzene-d5	6.63	82	1003346	72.64	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	321251	112.95	ng	0.00
44) 2-Fluorobiphenyl	8.82	172	1751762	74.24	ng	0.00
78) Terphenyl-d14	13.44	244	1359903	70.12	ng	0.00
Target Compounds						Qvalue
10) Phenol	5.58	94	67823	4.15	ng	89
49) Dimethylphthalate	9.30	163	153530	6.00	ng	98
51) Acenaphthene	9.67	154	78847	3.58	ng	99
57) Fluorene	10.30	166	73877m	2.97	ng	
70) Phenanthrene	11.49	178	1163268	33.75	ng	99
71) Anthracene	11.54	178	246294	6.94	ng	99
74) Fluoranthene	12.96	202	1133122	32.10	ng	98
77) Pyrene	13.23	202	1291623	40.94	ng	99
80) Benzo(a)anthracene	14.70	228	466649	18.41	ng	99
82) Chrysene	14.74	228	462439	19.66	ng	97
85) Indeno(1,2,3-cd)pyrene	17.73	276	125681	6.17	ng	97
87) Benzo(b)fluoranthene	15.93	252	346417m	17.58	ng	
88) Benzo(k)fluoranthene	15.96	252	92858m	4.81	ng	
89) Benzo(a)pyrene	16.28	252	261101	14.38	ng	97
91) Benzo(a,h,i)perylene	18.14	276	133453	8.61	ng	98

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

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