

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012717\
 Data File : BF092595.D
 Acq On : 28 Jan 2017 00:10
 Operator : SJ/MA
 Sample : I1531-01
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EDC-HPRR-S1C1

Quant Time: Jan 28 04:41:37 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 24 13:48:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	176604m	20.00	ng	0.00
21) Naphthalene-d8	8.27	136	713787	20.00	ng	0.00
38) Acenaphthene-d10	10.03	164	327493	20.00	ng	0.00
63) Phenanthrene-d10	11.53	188	460223	20.00	ng	0.00
75) Chrysene-d12	14.17	240	360695	20.00	ng	0.00
86) Perylene-d12	15.65	264	281855	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	1713654	150.97	ng	0.06
7) Phenol-d6	6.62	99	1871476	129.44	ng	0.03
23) Nitrobenzene-d5	7.55	82	1359429	104.04	ng	0.01
41) 2,4,6-Tribromophenol	10.83	330	315409	141.33	ng	0.01
44) 2-Fluorobiphenyl	9.36	172	2122651	119.85	ng	0.00
78) Terphenyl-d14	13.12	244	1377461	101.69	ng	0.00
Target Compounds						
10) Phenol	6.64	94	82051	4.70	ng	86
49) Dimethylphthalate	9.74	163	272289	13.07	ng	99
70) Phenanthrene	11.55	178	429500	16.95	ng	98
71) Anthracene	11.60	178	117266	4.73	ng	97
74) Fluoranthene	12.74	202	620316	25.02	ng	96
77) Pyrene	12.97	202	563298	21.55	ng	99
80) Benzo(a)anthracene	14.16	228	245733m	12.68	ng	
82) Chrysene	14.19	228	241647m	11.67	ng	
85) Indeno(1,2,3-cd)pyrene	17.15	276	77269	5.04	ng	93
87) Benzo(b)fluoranthene	15.22	252	227024m	12.55	ng	
88) Benzo(k)fluoranthene	15.23	252	89395m	5.94	ng	
89) Benzo(a)pyrene	15.59	252	163538	10.68	ng	# 93
91) Benzo(g,h,i)perylene	17.60	276	72839	5.82	ng	# 89

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

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