

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042617\
 Data File : BF094657.D
 Acq On : 26 Apr 2017 19:56
 Operator : SJ/MA
 Sample : I2813-04
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-W9-SSW

Quant Time: Apr 27 05:44:16 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 17 14:59:23 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.75	152	134837	20.00	ng	0.00
21) Naphthalene-d8	7.25	136	557308	20.00	ng	-0.01
38) Acenaphthene-d10	9.38	164	232600	20.00	ng	-0.02
63) Phenanthrene-d10	11.20	188	398076	20.00	ng	-0.02
75) Chrysene-d12	14.45	240	270467	20.00	ng	-0.01
86) Perylene-d12	16.08	264	250721	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.30	112	916049	112.71	ng	0.00
7) Phenol-d6	5.38	99	1154127	120.46	ng	0.00
23) Nitrobenzene-d5	6.41	82	701481	77.72	ng	0.00
41) 2,4,6-Tribromophenol	10.36	330	205998	113.23	ng	-0.01
44) 2-Fluorobiphenyl	8.57	172	1174477	74.29	ng	-0.01
78) Terphenyl-d14	13.18	244	644270	63.67	ng	-0.01
Target Compounds						Qvalue
10) Phenol	5.40	94	65653	5.58	ng	87
49) Dimethylphthalate	9.07	163	161459	9.31	ng	98
51) Acenaphthene	9.42	154	30433	2.16	ng	99
57) Fluorene	10.05	166	51933	3.25	ng	97
70) Phenanthrene	11.24	178	723545	32.28	ng	98
71) Anthracene	11.29	178	210458	9.08	ng	96
74) Fluoranthene	12.69	202	656473	28.26	ng	99
77) Pyrene	12.97	202	601547	31.83	ng	98
80) Benzo(a)anthracene	14.44	228	202830	12.90	ng	99
82) Chrysene	14.48	228	179078	12.46	ng	98
85) Indeno(1,2,3-cd)pyrene	17.35	276	82138	6.01	ng	95
87) Benzo(b)fluoranthene	15.68	252	178134m	11.61	ng	
88) Benzo(k)fluoranthene	15.71	252	71423m	4.92	ng	
89) Benzo(a)pyrene	16.02	252	160106	11.22	ng	97
91) Benzo(a,h,i)perylene	17.73	276	76050	6.30	ng	95

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

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