

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041917\
 Data File : BF094526.D
 Acq On : 19 Apr 2017 20:03
 Operator : SJ/MA
 Sample : I2750-04
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-T5-WSW

Quant Time: Apr 20 07:58:02 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	5.77	152	125345	20.00	ng	0.00
21) Naphthalene-d8	7.26	136	517873	20.00	ng	0.00
38) Acenaphthene-d10	9.40	164	206942	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	413916	20.00	ng	0.00
75) Chrysene-d12	14.47	240	274966	20.00	ng	0.00
86) Perylene-d12	16.09	264	251705	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	4.32	112	1058169	140.06	ng	0.02
7) Phenol-d6	5.40	99	1401679	157.37	ng	0.00
23) Nitrobenzene-d5	6.42	82	849396	101.28	ng	0.00
41) 2,4,6-Tribromophenol	10.38	330	280218	173.12	ng	0.00
44) 2-Fluorobiphenyl	8.59	172	1375251	97.78	ng	0.00
78) Terphenyl-d14	13.19	244	1024244	99.57	ng	0.00

Target Compounds					Qvalue
10) Phenol	5.41	94	31966	2.92 ng	90
31) Naphthalene	7.29	128	219231	8.81 ng	97
37) 2-Methylnaphthalene	8.12	142	73247	4.30 ng	99
49) Dimethylphthalate	9.08	163	414464	26.87 ng	99
51) Acenaphthene	9.43	154	122006	9.72 ng	98
54) Dibenzofuran	9.65	168	144650	7.93 ng	97
57) Fluorene	10.07	166	175393	12.35 ng	100
70) Phenanthrene	11.25	178	1328035	56.98 ng	98
71) Anthracene	11.31	178	295975	12.28 ng	97
72) Carbazole	11.51	167	167538	7.40 ng	97
74) Fluoranthene	12.71	202	1135715	47.01 ng	99
77) Pyrene	12.98	202	1065819	55.48 ng	98
80) Benzo(a)anthracene	14.45	228	321535	20.12 ng	98
82) Chrysene	14.49	228	316506	21.67 ng	97
85) Indeno(1,2,3-cd)pyrene	17.36	276	187919	13.52 ng	93
87) Benzo(b)fluoranthene	15.69	252	340272m	22.10 ng	
88) Benzo(k)fluoranthene	15.71	252	140431m	9.64 ng	
89) Benzo(a)pyrene	16.03	252	303075	21.16 ng	97
90) Dibenzo(a,h)anthracene	17.37	278	40806m	3.43 ng	
91) Benzo(g,h,i)perylene	17.73	276	199803	16.50 ng	99

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

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