

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072117\
 Data File : BF096940.D
 Acq On : 21 Jul 2017 13:22
 Operator : SJ/JU
 Sample : I4253-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-O-P-6-7-SSW

Quant Time: Jul 24 11:18:25 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF071317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 18 13:33:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.55	152	114594	20.00	ng	-0.05
21) Naphthalene-d8	7.83	136	454873	20.00	ng	-0.05
38) Acenaphthene-d10	9.58	164	210643	20.00	ng	-0.05
63) Phenanthrene-d10	11.06	188	289980	20.00	ng	-0.05
75) Chrysene-d12	13.69	240	208332	20.00	ng	-0.05
86) Perylene-d12	15.05	264	203010	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	872282	114.45	ng	-0.03
7) Phenol-d6	6.22	99	1103687	120.68	ng	-0.04
23) Nitrobenzene-d5	7.12	82	636427	86.66	ng	-0.05
41) 2,4,6-Tribromophenol	10.37	330	211834	117.82	ng	-0.05
44) 2-Fluorobiphenyl	8.92	172	1058593	79.40	ng	-0.05
78) Terphenyl-d14	12.65	244	680343	56.64	ng	-0.05
Target Compounds						Qvalue
10) Phenol	6.23	94	22870	2.212	ng	91
49) Dimethylphthalate	9.31	163	172253	10.860	ng	99
70) Phenanthrene	11.08	178	308662	19.576	ng	98
71) Anthracene	11.13	178	84803	5.320	ng	98
72) Carbazole	11.29	167	37121	2.405	ng	97
74) Fluoranthene	12.27	202	434168	28.768	ng	99
77) Pyrene	12.49	202	441468	23.349	ng	98
80) Benzo(a)anthracene	13.67	228	280758	21.629	ng	97
82) Chrysene	13.71	228	238723m	18.681	ng	
85) Indeno(1,2,3-cd)pyrene	16.32	276	202201	19.402	ng	94
87) Benzo(b)fluoranthene	14.69	252	384644m	31.415	ng	
88) Benzo(k)fluoranthene	14.70	252	114943m	9.914	ng	
89) Benzo(a)pyrene	15.00	252	308767	27.495	ng	95
90) Dibenzo(a,h)anthracene	16.32	278	49987m	4.838	ng	
91) Benzo(a,h,i)perylene	16.69	276	194328	17.562	ng	97

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

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