

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030617\
 Data File : BF093407.D
 Acq On : 6 Mar 2017 18:13
 Operator : SJ/MA
 Sample : I2088-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-004-B

Quant Time: Mar 07 00:24:50 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 07 00:03:56 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	196386	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	769571	20.00	ng	0.00
38) Acenaphthene-d10	9.81	164	339565	20.00	ng	0.00
63) Phenanthrene-d10	11.29	188	540357	20.00	ng	0.00
75) Chrysene-d12	13.93	240	408833	20.00	ng	0.00
86) Perylene-d12	15.35	264	390234	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.37	112	1580531	138.07	ng	0.02
7) Phenol-d6	6.44	99	2216700	150.50	ng	0.00
23) Nitrobenzene-d5	7.35	82	1413962	102.32	ng	0.00
41) 2,4,6-Tribromophenol	10.61	330	309446	126.00	ng	0.00
44) 2-Fluorobiphenyl	9.13	172	1795815	95.81	ng	0.00
78) Terphenyl-d14	12.88	244	1302161	74.84	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.45	94	134202	7.79	ng	# 66
49) Dimethylphthalate	9.53	163	172015	7.52	ng	99
70) Phenanthrene	11.32	178	227050	7.33	ng	97
71) Anthracene	11.37	178	71281	2.29	ng	95
74) Fluoranthene	12.50	202	382110	11.97	ng	98
77) Pyrene	12.73	202	397525	12.99	ng	98
80) Benzo(a)anthracene	13.92	228	177749m	7.11	ng	
82) Chrysene	13.96	228	168308m	7.19	ng	
85) Indeno(1,2,3-cd)pyrene	16.72	276	88284	4.87	ng	99
87) Benzo(b)fluoranthene	14.95	252	201438m	7.84	ng	
88) Benzo(k)fluoranthene	14.97	252	57014m	2.57	ng	
89) Benzo(a)pyrene	15.29	252	153155	6.89	ng	# 97
91) Benzo(a,h,i)perylene	17.13	276	79083	4.36	ng	# 90

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

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