

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021717\
 Data File : BF093023.D
 Acq On : 17 Feb 2017 21:04
 Operator : SJ/MA
 Sample : I1892-05
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP22-COMP16

Manual Integrations
 APPROVED

mohammad
 2/20/2017 1:00:01 PM

Quant Time: Feb 17 23:36:19 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 17 17:31:55 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	164321	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	654731	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	299505	20.00	ng	0.00
63) Phenanthrene-d10	11.38	188	536823	20.00	ng	0.00
75) Chrysene-d12	14.01	240	331493	20.00	ng	-0.01
86) Perylene-d12	15.44	264	313409	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	1173918	122.57	ng	0.01
7) Phenol-d6	6.50	99	1351954	109.70	ng	0.01
23) Nitrobenzene-d5	7.42	82	809554	68.86	ng	0.00
41) 2,4,6-Tribromophenol	10.68	330	263768	121.77	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	1368092	82.75	ng	0.00
78) Terphenyl-d14	12.97	244	1164448	82.54	ng	0.00
Target Compounds						
10) Phenol	6.51	94	144292	10.01	ng	# 69
49) Dimethylphthalate	9.61	163	141638	7.02	ng	# 98
70) Phenanthrene	11.40	178	151614	4.93	ng	95
74) Fluoranthene	12.59	202	164768	5.19	ng	92
77) Pyrene	12.82	202	157871	6.36	ng	97
80) Benzo(a)anthracene	14.00	228	100171	4.94	ng	99
82) Chrysene	14.03	228	71117	3.75	ng	96
85) Indeno(1,2,3-cd)pyrene	16.84	276	70604	4.80	ng	99
87) Benzo(b)fluoranthene	15.03	252	131691m	6.38	ng	
89) Benzo(a)pyrene	15.38	252	110290	6.18	ng	98
91) Benzo(g,h,i)perylene	17.27	276	72324	4.96	ng	# 90

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

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