

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071217\
 Data File : BF096699.D
 Acq On : 12 Jul 2017 18:43
 Operator : SJ/JU
 Sample : I4144-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP1-0-3

Manual Integrations
 APPROVED

mohammad
 7/13/2017 1:37:55 PM

Quant Time: Jul 13 07:53:33 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 07 13:02:33 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	123228	20.00	ng	-0.03
21) Naphthalene-d8	7.95	136	477215	20.00	ng	-0.04
38) Acenaphthene-d10	9.70	164	193504	20.00	ng	-0.04
63) Phenanthrene-d10	11.18	188	319186	20.00	ng	-0.03
75) Chrysene-d12	13.81	240	197050	20.00	ng	-0.04
86) Perylene-d12	15.20	264	172283	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.28	112	745078	89.24	ng	-0.03
7) Phenol-d6	6.31	99	899342	87.62	ng	-0.04
23) Nitrobenzene-d5	7.23	82	565777	71.24	ng	-0.04
41) 2,4,6-Tribromophenol	10.49	330	145121	96.01	ng	-0.04
44) 2-Fluorobiphenyl	9.03	172	869937	76.87	ng	-0.04
78) Terphenyl-d14	12.76	244	517124	56.08	ng	-0.04
Target Compounds						Qvalue
31) Naphthalene	7.97	128	67706	3.005	ng	99
37) 2-Methylnaphthalene	8.66	142	53970	3.763	ng	95
48) Acenaphthylene	9.56	152	46333	2.367	ng	95
49) Dimethylphthalate	9.42	163	291530	20.182	ng	100
51) Acenaphthene	9.73	154	31767	2.632	ng	97
54) Dibenzofuran	9.90	168	38997	2.447	ng	# 94
57) Fluorene	10.24	166	29663	2.636	ng	# 95
70) Phenanthrene	11.20	178	417666	24.206	ng	99
71) Anthracene	11.25	178	115822	6.587	ng	97
72) Carbazole	11.40	167	46408	2.624	ng	99
74) Fluoranthene	12.39	202	445225	26.318	ng	98
77) Pyrene	12.62	202	371937	23.516	ng	99
80) Benzo(a)anthracene	13.80	228	193524	16.960	ng	94
82) Chrysene	13.83	228	187654	15.704	ng	96
85) Indeno(1,2,3-cd)pyrene	16.53	276	85707	9.087	ng	94
87) Benzo(b)fluoranthene	14.82	252	243156m	22.524	ng	
88) Benzo(k)fluoranthene	14.84	252	89923m	9.312	ng	
89) Benzo(a)pyrene	15.14	252	132602	13.758	ng	99
90) Dibenzo(a,h)anthracene	16.53	278	23458m	3.094	ng	
91) Benzo(g,h,i)perylene	16.93	276	74852	9.527	ng	97

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

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