

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090617\
 Data File : BF098290.D
 Acq On : 6 Sep 2017 23:26
 Operator : SJ/JU
 Sample : I5053-06
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G-62A-1.5-3

Manual Integrations
 APPROVED

Sohil
 9/7/2017 5:44:39 PM

Quant Time: Sep 07 06:48:44 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 05 17:28:20 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	101659	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	380445	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	148235	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	235783	20.00	ng	0.00
75) Chrysene-d12	13.85	240	218272	20.00	ng	-0.01
86) Perylene-d12	15.26	264	155102	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	683850	102.32	ng	0.00
7) Phenol-d6	6.35	99	896577	98.73	ng	0.00
23) Nitrobenzene-d5	7.26	82	536046	76.54	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	120891	93.99	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	748313	74.17	ng	-0.01
78) Terphenyl-d14	12.80	244	496212	47.41	ng	-0.01
Target Compounds						Qvalue
37) 2-Methylnaphthalene	8.69	142	24891	2.056	ng	98
49) Dimethylphthalate	9.45	163	129106	11.915	ng	98
70) Phenanthrene	11.24	178	110690	8.810	ng	97
71) Anthracene	11.29	178	30661	2.406	ng	99
74) Fluoranthene	12.42	202	158923	12.118	ng	98
77) Pyrene	12.65	202	145002	8.122	ng	97
80) Benzo(a)anthracene	13.84	228	77915	5.956	ng	98
82) Chrysene	13.87	228	94553	7.266	ng	97
85) Indeno(1,2,3-cd)pyrene	16.62	276	27877	2.912	ng	93
87) Benzo(b)fluoranthene	14.86	252	95953m	9.815	ng	
88) Benzo(k)fluoranthene	14.88	252	34040m	3.706	ng	
89) Benzo(a)pyrene	15.20	252	46883	5.417	ng	# 90
91) Benzo(a,h,i)perylene	17.03	276	28038	3.814	ng	91

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

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