

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091417\
 Data File : BF098536.D
 Acq On : 14 Sep 2017 19:02
 Operator : SJ/JU
 Sample : I5160-09
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G-8B-2

Manual Integrations
 APPROVED

Sohil
 9/15/2017 5:43:49 PM

Quant Time: Sep 15 05:37:18 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF091117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 02:14:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	135783	20.00	ng	-0.02
21) Naphthalene-d8	7.93	136	569171	20.00	ng	-0.02
38) Acenaphthene-d10	9.68	164	265966	20.00	ng	-0.02
63) Phenanthrene-d10	11.16	188	505484	20.00	ng	-0.02
75) Chrysene-d12	13.79	240	311797	20.00	ng	-0.02
86) Perylene-d12	15.19	264	294937	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	1378832	150.14	ng	-0.01
7) Phenol-d6	6.30	99	1695494	145.41	ng	-0.02
23) Nitrobenzene-d5	7.22	82	1071069	104.91	ng	-0.02
41) 2,4,6-Tribromophenol	10.47	330	352252	198.44	ng	-0.02
44) 2-Fluorobiphenyl	9.01	172	1808194	115.23	ng	-0.02
78) Terphenyl-d14	12.75	244	1542655	106.79	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.40	163	200156	9.722	ng	99
70) Phenanthrene	11.19	178	178069	6.645	ng	96
74) Fluoranthene	12.37	202	350911	13.081	ng	99
77) Pyrene	12.60	202	299312	12.169	ng	98
80) Benzo(a)anthracene	13.79	228	148750	8.137	ng	97
82) Chrysene	13.82	228	180370	9.761	ng	97
85) Indeno(1,2,3-cd)pyrene	16.52	276	113428	8.749	ng	92
87) Benzo(b)fluoranthene	14.80	252	230511m	12.430	ng	
88) Benzo(k)fluoranthene	14.82	252	65227m	3.768	ng	
89) Benzo(a)pyrene	15.13	252	137762	8.338	ng	97
90) Dibenzo(a,h)anthracene	16.53	278	30603	2.546	ng	# 86
91) Benzo(a,h,i)perylene	16.92	276	115507	9.555	ng	95

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

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