

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090817\
 Data File : BF098360.D
 Acq On : 8 Sep 2017 19:43
 Operator : SJ/JU
 Sample : I5071-18
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G57A-1.5-3

Manual Integrations
 APPROVED

Sohil
 9/11/2017 4:47:43 PM

Quant Time: Sep 11 11:39:16 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	6.68	152	98103	20.00	ng	-0.02
21) Naphthalene-d8	7.96	136	375457	20.00	ng	-0.02
38) Acenaphthene-d10	9.72	164	143255	20.00	ng	-0.02
63) Phenanthrene-d10	11.20	188	229372	20.00	ng	-0.02
75) Chrysene-d12	13.84	240	222565	20.00	ng	-0.02
86) Perylene-d12	15.24	264	177906	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	781574	121.18	ng	0.00
7) Phenol-d6	6.33	99	1064624	121.49	ng	-0.02
23) Nitrobenzene-d5	7.25	82	629618	91.10	ng	-0.02
41) 2,4,6-Tribromophenol	10.51	330	154412	124.23	ng	-0.02
44) 2-Fluorobiphenyl	9.04	172	948117	97.24	ng	-0.02
78) Terphenyl-d14	12.79	244	663273	62.15	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	7.99	128	43127	2.213	ng	98
37) 2-Methylnaphthalene	8.68	142	28814	2.411	ng	97
49) Dimethylphthalate	9.44	163	152854	14.597	ng	98
70) Phenanthrene	11.22	178	92309	7.552	ng	98
71) Anthracene	11.27	178	37424	3.019	ng	98
74) Fluoranthene	12.41	202	163788	12.839	ng	98
77) Pyrene	12.64	202	152810	8.395	ng	97
80) Benzo(a)anthracene	13.83	228	99238	7.440	ng	93
82) Chrysene	13.86	228	121049	9.122	ng	96
85) Indeno(1,2,3-cd)pyrene	16.60	276	44787m	4.588	ng	
87) Benzo(b)fluoranthene	14.85	252	157434m	14.039	ng	
88) Benzo(k)fluoranthene	14.87	252	50038m	4.749	ng	
89) Benzo(a)pyrene	15.19	252	73059	7.359	ng	# 95
91) Benzo(a,h,i)perylene	17.00	276	36642	4.345	ng	92

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

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