

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091217\
 Data File : BF098453.D
 Acq On : 12 Sep 2017 16:35
 Operator : SJ/JU
 Sample : I5119-01RE
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SASP2P-TP-116A(0-5)COMPRES

Manual Integrations
 APPROVED

Sohil
 9/13/2017 5:00:03 PM

Quant Time: Sep 13 01:28:46 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	6.66	152	228022	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	927395	20.00	ng	0.00
38) Acenaphthene-d10	9.69	164	384391	20.00	ng	-0.01
63) Phenanthrene-d10	11.17	188	597535	20.00	ng	-0.01
75) Chrysene-d12	13.81	240	426160	20.00	ng	0.00
86) Perylene-d12	15.21	264	363115	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	520102	33.72	ng	-0.01
7) Phenol-d6	6.32	99	766497	39.14	ng	0.00
23) Nitrobenzene-d5	7.23	82	407481	24.50	ng	-0.01
41) 2,4,6-Tribromophenol	10.49	330	285744	111.38	ng	0.00
44) 2-Fluorobiphenyl	9.02	172	1023571	45.13	ng	-0.01
78) Terphenyl-d14	12.76	244	1315455	66.63	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.42	163	165832	5.573	ng	99
70) Phenanthrene	11.20	178	486620	15.362	ng	98
71) Anthracene	11.25	178	134936	4.149	ng	98
74) Fluoranthene	12.39	202	1312411	41.386	ng	98
77) Pyrene	12.62	202	1332138	39.625	ng	99
80) Benzo(a)anthracene	13.80	228	743312	29.750	ng	98
82) Chrysene	13.84	228	649351	25.711	ng	97
85) Indeno(1,2,3-cd)pyrene	16.55	276	276167	15.586	ng	95
87) Benzo(b)fluoranthene	14.82	252	769770m	33.715	ng	
88) Benzo(k)fluoranthene	14.84	252	289221m	13.569	ng	
89) Benzo(a)pyrene	15.16	252	582307	28.628	ng	98
90) Dibenzo(a,h)anthracene	16.56	278	71769m	4.850	ng	
91) Benzo(a,h,i)perylene	16.96	276	272028	18.278	ng	99

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

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