

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110117\
 Data File : BF100066.D
 Acq On : 2 Nov 2017 1:00
 Operator : SJ/JU
 Sample : I6079-14 5X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Q3-COMP

Manual Integrations
 APPROVED

Sohil
 11/2/2017 4:29:55 PM

Quant Time: Nov 02 05:10:54 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 18:13:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	68807	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	280527	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	119783	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	188829	20.00	ng	0.01
75) Chrysene-d12	14.32	240	133837m	20.00	ng	0.35
86) Perylene-d12	16.22	264	120782m	20.00	ng	0.75
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	123144	28.10	ng	0.01
7) Phenol-d6	6.43	99	156383	30.09	ng	0.00
23) Nitrobenzene-d5	7.35	82	84257	19.34	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	29295	26.84	ng	0.00
44) 2-Fluorobiphenyl	9.13	172	186501	24.02	ng	0.00
78) Terphenyl-d14	13.02	244	118573	19.36	ng	0.12
Target Compounds						
49) Dimethylphthalate	9.54	163	49053	5.171	ng	99
70) Phenanthrene	11.35	178	105062	9.859	ng	98
71) Anthracene	11.40	178	27428	2.536	ng	97
74) Fluoranthene	12.60	202	121883	11.006	ng	99
77) Pyrene	12.86	202	103841	10.204	ng	98
80) Benzo(a)anthracene	14.31	228	50745m	5.981	ng	
82) Chrysene	14.36	228	40843m	5.120	ng	
87) Benzo(b)fluoranthene	15.72	252	47108m	6.177	ng	
88) Benzo(k)fluoranthene	15.75	252	21330m	2.966	ng	
91) Benzo(q,h,i)perylene	17.81	276	17113m	2.873	ng	

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

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