

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF021918\
 Data File : BF103082.D
 Acq On : 19 Feb 2018 15:12
 Operator : SJ/JU
 Sample : J1604-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3

Manual Integrations
 APPROVED

Sohil
 2/20/2018 10:20:28 AM

Quant Time: Feb 19 17:35:12 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 16 14:34:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	177143	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	729679	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	316855	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	479382	20.00	ng	0.00
75) Chrysene-d12	14.04	240	314935	20.00	ng	-0.01
86) Perylene-d12	15.54	264	336212	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1091719	93.59	ng	0.01
7) Phenol-d6	6.51	99	1314686	93.61	ng	0.00
23) Nitrobenzene-d5	7.44	82	855548	81.34	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	251032	98.43	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1272798	66.66	ng	0.00
78) Terphenyl-d14	12.99	244	859390	64.61	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.62	163	74033	2.997	ng	98
70) Phenanthrene	11.43	178	108034	4.046	ng	97
74) Fluoranthene	12.62	202	184168	6.856	ng	96
77) Pyrene	12.85	202	147701	5.981	ng	97
80) Benzo(a)anthracene	14.04	228	64977	3.281	ng	98
82) Chrysene	14.07	228	76203	3.817	ng	96
87) Benzo(b)fluoranthene	15.10	252	100034m	4.589	ng	
89) Benzo(a)pyrene	15.47	252	57491	2.930	ng	# 96

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

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