

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF120117\
 Data File : BF101072.D
 Acq On : 1 Dec 2017 19:18
 Operator : SJ/JU
 Sample : I6622-10
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DBNW1A-2-4

Manual Integrations
 APPROVED

Sohil
 12/4/2017 11:56:48 AM

Quant Time: Dec 01 23:43:02 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 01 15:52:19 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	54009	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	207555	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	89125	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	141805	20.00	ng	0.00
75) Chrysene-d12	14.02	240	122872	20.00	ng	0.00
86) Perylene-d12	15.49	264	103638	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	403633	123.49	ng	0.01
7) Phenol-d6	6.49	99	487513	122.86	ng	0.00
23) Nitrobenzene-d5	7.42	82	295494	84.93	ng	0.00
41) 2,4,6-Tribromophenol	10.67	330	94244	88.03	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	513903	78.89	ng	0.00
78) Terphenyl-d14	12.96	244	376122	66.95	ng	0.00
Target Compounds						
10) Phenol	6.50	94	16711	3.787	ng	96
49) Dimethylphthalate	9.60	163	50780	7.065	ng	99
70) Phenanthrene	11.40	178	17815	2.281	ng	98
74) Fluoranthene	12.59	202	26784	3.094	ng	95
77) Pyrene	12.82	202	24427	2.881	ng	96
80) Benzo(a)anthracene	14.00	228	15924	2.053	ng	91
82) Chrysene	14.04	228	16158m	2.243	ng	
87) Benzo(b)fluoranthene	15.06	252	19025m	3.065	ng	
89) Benzo(a)pyrene	15.43	252	11904	2.109	ng	96

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

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