

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112817\
 Data File : BF100958.D
 Acq On : 29 Nov 2017 3:03
 Operator : SJ/JU
 Sample : I6579-11
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HAR-2-E(7-8)

Manual Integrations
 APPROVED

Sohil
 11/29/2017 4:29:25 PM

Quant Time: Nov 29 08:10:07 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 28 15:21:02 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	66448	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	243662	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	98988	20.00	ng	0.00
63) Phenanthrene-d10	11.38	188	164807	20.00	ng	0.00
75) Chrysene-d12	14.02	240	155421	20.00	ng	0.00
86) Perylene-d12	15.50	264	118882	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	488015	117.73	ng	0.01
7) Phenol-d6	6.49	99	589024	115.23	ng	0.00
23) Nitrobenzene-d5	7.42	82	338789	91.92	ng	0.00
41) 2,4,6-Tribromophenol	10.69	330	122826	121.77	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	524531	80.69	ng	0.00
78) Terphenyl-d14	12.97	244	404319	59.77	ng	0.00
Target Compounds						
10) Phenol	6.50	94	15225	2.837	ng	89
49) Dimethylphthalate	9.60	163	35800	4.737	ng	98
70) Phenanthrene	11.40	178	40161	4.628	ng	99
74) Fluoranthene	12.59	202	53013	5.765	ng	95
77) Pyrene	12.82	202	43419	3.919	ng	98
80) Benzo(a)anthracene	14.01	228	27759	2.966	ng	100
82) Chrysene	14.05	228	25116	2.771	ng	96
87) Benzo(b)fluoranthene	15.06	252	27397m	3.890	ng	
89) Benzo(a)pyrene	15.43	252	18884	3.024	ng	98

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

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