

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103118\
 Data File : BF110338.D
 Acq On : 31 Oct 2018 17:41
 Operator : JU/SJ
 Sample : J5728-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EG-DRUM

Quant Time: Nov 01 04:20:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 31 14:18:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	80223	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	320056	20.00	ng	0.00
39) Acenaphthene-d10	10.02	164	138683	20.00	ng	0.00
64) Phenanthrene-d10	11.52	188	240678	20.00	ng	0.00
76) Chrysene-d12	14.17	240	167606	20.00	ng	0.00
87) Perylene-d12	15.71	264	123900	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	664471	134.88	ng	0.01
7) Phenol-d6	6.60	99	825280	130.97	ng	0.00
23) Nitrobenzene-d5	7.54	82	491921	91.16	ng	0.00
42) 2,4,6-Tribromophenol	10.82	330	173900	126.59	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	878896	99.51	ng	0.00
79) Terphenyl-d14	13.10	244	699055	86.74	ng	0.00
Target Compounds						
10) Phenol	6.62	94	55313	8.212	ng	# 64
50) Dimethylphthalate	9.72	163	134971	13.184	ng	# 98
71) Phenanthrene	11.54	178	94275	7.486	ng	99
75) Fluoranthene	12.73	202	41863	3.275	ng	99
78) Pyrene	12.96	202	73006	6.076	ng	97
81) Benzo(a)anthracene	14.16	228	23965	2.411	ng	# 86
83) Chrysene	14.19	228	23055	2.442	ng	# 96
90) Benzo(a)pyrene	15.64	252	13881	2.108	ng	# 91

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

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