

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112718\
 Data File : BF111012.D
 Acq On : 27 Nov 2018 16:24
 Operator : JU/SJ
 Sample : J6060-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20180945-COMPOSITE-4

Manual Integrations
 APPROVED

mohammad
 11/28/2018 4:42:16 PM

Quant Time: Nov 27 17:37:50 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 27 15:32:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	44692	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	161536	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	65184	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	126304	20.00	ng	0.00
76) Chrysene-d12	14.14	240	102483	20.00	ng	0.00
87) Perylene-d12	15.67	264	67306	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	347871	129.08	ng	0.01
7) Phenol-d6	6.57	99	405514	115.46	ng	0.00
23) Nitrobenzene-d5	7.50	82	243395	86.23	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	75271	116.27	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	368612	81.93	ng	0.00
79) Terphenyl-d14	13.06	244	348422	66.18	ng	0.00
Target Compounds						
10) Phenol	6.58	94	21686	5.483	ng	87
50) Dimethylphthalate	9.69	163	50506	10.565	ng	99
71) Phenanthrene	11.50	178	26635	3.982	ng	98
75) Fluoranthene	12.70	202	33156	4.822	ng	98
78) Pyrene	12.93	202	37530	5.014	ng	98
81) Benzo(a)anthracene	14.13	228	13662m	2.258	ng	
83) Chrysene	14.16	228	13475	2.262	ng	99
88) Benzo(b)fluoranthene	15.22	252	8340m	2.109	ng	
90) Benzo(a)pyrene	15.62	252	7527	2.045	ng	# 74

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

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