

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111317\
 Data File : BF100503.D
 Acq On : 13 Nov 2017 15:56
 Operator : SJ/JU
 Sample : I6409-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-110917-02-FREEZE-PIT

Manual Integrations
 APPROVED

SOHIL
 11/14/2017 5:20:18 PM

Quant Time: Nov 14 01:16:09 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 13 14:08:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	92958	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	354792	20.00	ng	0.00
38) Acenaphthene-d10	9.93	164	156540	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	260947	20.00	ng	0.00
75) Chrysene-d12	14.05	240	191769	20.00	ng	0.00
86) Perylene-d12	15.53	264	162542	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	502774	86.70	ng	0.00
7) Phenol-d6	6.52	99	620685	86.80	ng	0.00
23) Nitrobenzene-d5	7.45	82	391854	73.02	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	148757	93.26	ng	0.00
44) 2-Fluorobiphenyl	9.25	172	663392	64.53	ng	0.00
78) Terphenyl-d14	13.00	244	410024	49.13	ng	0.00
Target Compounds						
10) Phenol	6.53	94	30820	4.105	ng	97
31) Naphthalene	8.19	128	224948	13.164	ng	99
49) Dimethylphthalate	9.63	163	73258	6.129	ng	99
74) Fluoranthene	12.62	202	36250	2.490	ng	96
77) Pyrene	12.85	202	43224	3.162	ng	99
82) Chrysene	14.07	228	22785	2.037	ng	# 91
87) Benzo(b)fluoranthene	15.09	252	26343m	2.736	ng	
89) Benzo(a)pyrene	15.46	252	20931	2.452	ng	# 92
91) Benzo(g,h,i)perylene	17.46	276	16324	2.389	ng	# 87

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

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