

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111317\
 Data File : BF100505.D
 Acq On : 13 Nov 2017 16:50
 Operator : SJ/JU
 Sample : I6301-05 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-110917-C

Manual Integrations
 APPROVED

SOHIL
 11/14/2017 5:20:23 PM

Quant Time: Nov 14 01:37:37 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	89975	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	332972	20.00	ng	0.00
38) Acenaphthene-d10	9.93	164	134064	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	216723	20.00	ng	0.00
75) Chrysene-d12	14.05	240	197724	20.00	ng	0.00
86) Perylene-d12	15.53	264	158041	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	251616	44.83	ng	0.00
7) Phenol-d6	6.51	99	303303	43.82	ng	0.00
23) Nitrobenzene-d5	7.45	82	182957	36.33	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	65937	48.27	ng	0.00
44) 2-Fluorobiphenyl	9.25	172	357928	40.65	ng	0.00
78) Terphenyl-d14	12.99	244	249856	29.04	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.63	163	50899	4.973	ng	Qvalue # 99
70) Phenanthrene	11.43	178	38221	3.350	ng	99
74) Fluoranthene	12.62	202	53953	4.461	ng	96
77) Pyrene	12.85	202	48916	3.471	ng	98
80) Benzo(a)anthracene	14.04	228	28013	2.353	ng	91
82) Chrysene	14.07	228	28475m	2.470	ng	
87) Benzo(b)fluoranthene	15.09	252	30874m	3.297	ng	
89) Benzo(a)pyrene	15.46	252	19908	2.398	ng	# 90

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

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