

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111117\
 Data File : BF100437.D
 Acq On : 11 Nov 2017 13:42
 Operator : SJ/JU
 Sample : I6299-03 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-01-110917

Manual Integrations
 APPROVED

SOHIL
 11/13/2017 4:54:16 PM

Quant Time: Nov 13 03:05:18 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 10 15:44:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	137044	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	481023	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	185446	20.00	ng	0.00
63) Phenanthrene-d10	11.43	188	310918	20.00	ng	0.00
75) Chrysene-d12	14.07	240	291882	20.00	ng	0.00
86) Perylene-d12	15.56	264	206499	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	158600	18.55	ng	0.00
7) Phenol-d6	6.52	99	184166	17.47	ng	-0.01
23) Nitrobenzene-d5	7.46	82	98327	13.51	ng	-0.01
41) 2,4,6-Tribromophenol	10.73	330	29659	15.70	ng	0.00
44) 2-Fluorobiphenyl	9.26	172	193248	15.87	ng	0.00
78) Terphenyl-d14	13.01	244	156793	12.34	ng	0.00
Target Compounds						
70) Phenanthrene	11.45	178	77645	4.743	ng	98
74) Fluoranthene	12.64	202	100379	5.786	ng	98
77) Pyrene	12.87	202	96492	4.638	ng	98
80) Benzo(a)anthracene	14.06	228	42731	2.431	ng	95
82) Chrysene	14.09	228	38634	2.270	ng	98
87) Benzo(b)fluoranthene	15.12	252	37934m	3.101	ng	
89) Benzo(a)pyrene	15.49	252	27376	2.524	ng	# 81

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

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