

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111117\
 Data File : BF100438.D
 Acq On : 11 Nov 2017 14:09
 Operator : SJ/JU
 Sample : I6407-01 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 3091-4235

Manual Integrations
 APPROVED

SOHIL
 11/13/2017 4:54:18 PM

Quant Time: Nov 13 03:09:45 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	160267	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	601357	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	249584	20.00	ng	0.00
63) Phenanthrene-d10	11.43	188	403749	20.00	ng	0.00
75) Chrysene-d12	14.07	240	334133	20.00	ng	0.00
86) Perylene-d12	15.55	264	249880	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	644523	64.47	ng	0.00
7) Phenol-d6	6.53	99	794328	64.43	ng	0.00
23) Nitrobenzene-d5	7.46	82	472503	51.95	ng	0.00
41) 2,4,6-Tribromophenol	10.73	330	157284	61.84	ng	0.00
44) 2-Fluorobiphenyl	9.26	172	875140	53.39	ng	0.00
78) Terphenyl-d14	13.01	244	649524	44.67	ng	0.00
Target Compounds						
10) Phenol	6.54	94	36701	2.836	ng	94
49) Dimethylphthalate	9.65	163	55408	2.908	ng	# 98
74) Fluoranthene	12.64	202	51325	2.278	ng	96
77) Pyrene	12.87	202	50959	2.139	ng	98
85) Indeno(1,2,3-cd)pyrene	17.04	276	48076	3.050	ng	94
87) Benzo(b)fluoranthene	15.12	252	37023m	2.501	ng	
89) Benzo(a)pyrene	15.49	252	27869	2.123	ng	# 85
91) Benzo(g,h,i)perylene	17.50	276	100658	9.580	ng	95

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

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