

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110217\
 Data File : BF100094.D
 Acq On : 2 Nov 2017 19:28
 Operator : SJ/JU
 Sample : I6249-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-1

Manual Integrations
 APPROVED

Sohil
 11/3/2017 11:19:40 AM

Quant Time: Nov 02 20:29:27 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 18:13:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	81903	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	331908	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	150859	20.00	ng	-0.01
63) Phenanthrene-d10	11.30	188	232359	20.00	ng	-0.01
75) Chrysene-d12	13.95	240	151918	20.00	ng	-0.02
86) Perylene-d12	15.41	264	148981	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	493312	94.56	ng	0.00
7) Phenol-d6	6.42	99	586257	94.78	ng	-0.01
23) Nitrobenzene-d5	7.34	82	326704	63.37	ng	-0.01
41) 2,4,6-Tribromophenol	10.60	330	118061	85.87	ng	-0.01
44) 2-Fluorobiphenyl	9.12	172	608794	62.26	ng	-0.01
78) Terphenyl-d14	12.88	244	391520	56.32	ng	-0.02
Target Compounds						
10) Phenol	6.43	94	16142	2.377	ng	Qvalue # 67
49) Dimethylphthalate	9.52	163	109935	9.202	ng	99
70) Phenanthrene	11.32	178	35085	2.676	ng	97
74) Fluoranthene	12.52	202	66045	4.846	ng	99
77) Pyrene	12.75	202	57616	4.988	ng	96
80) Benzo(a)anthracene	13.94	228	23888	2.480	ng	97
82) Chrysene	13.97	228	19988	2.208	ng	98
87) Benzo(b)fluoranthene	14.99	252	26646m	2.832	ng	
89) Benzo(a)pyrene	15.36	252	20285	2.400	ng	# 90

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

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