

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF101217\
 Data File : BF099414.D
 Acq On : 13 Oct 2017 00:15
 Operator : SJ/JU
 Sample : I5729-12 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 6E-(1-2)

Manual Integrations
 APPROVED

Sohil
 10/13/2017 3:06:34 PM

Quant Time: Oct 13 04:30:17 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 11 16:25:21 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	112134	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	433133	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	168564	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	239262	20.00	ng	0.00
75) Chrysene-d12	14.02	240	211754	20.00	ng	0.00
86) Perylene-d12	15.49	264	174997	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	154757	21.97	ng	0.00
7) Phenol-d6	6.48	99	187598	21.59	ng	0.00
23) Nitrobenzene-d5	7.41	82	104002	15.40	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	22697	16.46	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	211043	20.90	ng	0.00
78) Terphenyl-d14	12.95	244	131245	14.10	ng	0.00
Target Compounds						
48) Acenaphthylene	9.75	152	39881	2.395	ng	99
70) Phenanthrene	11.40	178	65139	5.086	ng	99
74) Fluoranthene	12.59	202	321352	24.779	ng	99
77) Pyrene	12.82	202	280990	17.402	ng	99
80) Benzo(a)anthracene	14.00	228	198479	15.479	ng	99
82) Chrysene	14.04	228	151464	12.408	ng	97
85) Indeno(1,2,3-cd)pyrene	16.96	276	64592	6.615	ng	99
87) Benzo(b)fluoranthene	15.06	252	226105m	21.014	ng	
88) Benzo(k)fluoranthene	15.08	252	65884m	6.200	ng	
89) Benzo(a)pyrene	15.42	252	133371	13.504	ng	# 95
90) Dibenzo(a,h)anthracene	16.97	278	20649m	2.612	ng	
91) Benzo(a,h,i)perylene	17.42	276	57606	7.373	ng	95

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

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