

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF101217\
 Data File : BF099436.D
 Acq On : 13 Oct 2017 12:02
 Operator : SJ/JU
 Sample : I5729-13DL 5X
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 7E-(1-2)DL

Manual Integrations
 APPROVED

Sohil
 10/15/2017 12:26:10 PM

Quant Time: Oct 13 14:58:41 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	99836	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	375925	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	134966	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	202362	20.00	ng	0.00
75) Chrysene-d12	14.02	240	183985	20.00	ng	0.00
86) Perylene-d12	15.49	264	142806	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	186138	29.68	ng	0.00
7) Phenol-d6	6.48	99	218201	28.20	ng	0.00
23) Nitrobenzene-d5	7.41	82	122000	20.81	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	24915	22.56	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	212129	26.24	ng	0.00
78) Terphenyl-d14	12.95	244	145796	18.02	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.60	163	27507	2.700	ng	96
70) Phenanthrene	11.40	178	216111	19.951	ng	99
71) Anthracene	11.45	178	35861	3.236	ng	100
72) Carbazole	11.61	167	22248	2.050	ng	99
74) Fluoranthene	12.59	202	369321	33.671	ng	97
77) Pyrene	12.82	202	321304	22.902	ng	98
80) Benzo(a)anthracene	14.00	228	138381	12.421	ng	99
82) Chrysene	14.04	228	131766	12.423	ng	98
85) Indeno(1,2,3-cd)pyrene	16.97	276	45585	5.373	ng	98
87) Benzo(b)fluoranthene	15.06	252	135157m	15.393	ng	
88) Benzo(k)fluoranthene	15.08	252	46606m	5.374	ng	
89) Benzo(a)pyrene	15.43	252	81781	10.147	ng	97
91) Benzo(a,h,i)perylene	17.42	276	42829	6.717	ng	93

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

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