

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF101217\
 Data File : BF099409.D
 Acq On : 12 Oct 2017 21:56
 Operator : SJ/JU
 Sample : I5729-09
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 5N-(2-3)

Manual Integrations
 APPROVED

Sohil
 10/13/2017 3:05:54 PM

Quant Time: Oct 13 03:02:26 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	119365	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	473118	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	183446	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	258954	20.00	ng	0.00
75) Chrysene-d12	14.02	240	222678	20.00	ng	0.00
86) Perylene-d12	15.49	264	192800	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	658388	87.81	ng	0.01
7) Phenol-d6	6.49	99	796766	86.14	ng	0.00
23) Nitrobenzene-d5	7.42	82	472894	64.10	ng	0.00
41) 2,4,6-Tribromophenol	10.68	330	111396	74.21	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	827653	75.32	ng	0.00
78) Terphenyl-d14	12.96	244	515998	52.70	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.60	163	110145	7.955	ng	99
70) Phenanthrene	11.40	178	282623	20.390	ng	97
71) Anthracene	11.45	178	49603	3.497	ng	99
74) Fluoranthene	12.59	202	527181	37.559	ng	97
77) Pyrene	12.82	202	504743	29.726	ng	98
80) Benzo(a)anthracene	14.01	228	388456	28.809	ng	99
82) Chrysene	14.04	228	427175	33.277	ng	98
85) Indeno(1,2,3-cd)pyrene	16.97	276	182948	17.816	ng	97
87) Benzo(b)fluoranthene	15.06	252	674110m	56.867	ng	
88) Benzo(k)fluoranthene	15.09	252	194724m	16.631	ng	
89) Benzo(a)pyrene	15.43	252	338196	31.081	ng	# 95
90) Dibenzo(a,h)anthracene	16.97	278	58138m	6.676	ng	
91) Benzo(a,h,i)perylene	17.42	276	160369	18.630	ng	97

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

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