

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042120\
 Data File : BF120102.D
 Acq On : 21 Apr 2020 15:13
 Operator : MA/SJ
 Sample : L2342-12
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-17-16-COMP

Manual Integrations
 APPROVED

mohammad
 4/24/2020 4:17:03 AM

Quant Time: Apr 21 15:56:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 21 13:48:22 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	121713	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	478874	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	244256	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	467467	20.00	ng	0.00
76) Chrysene-d12	13.86	240	297794	20.00	ng	-0.01
87) Perylene-d12	15.27	264	256586	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	623896	88.59	ng	0.00
7) Phenol-d6	6.33	99	848787	93.53	ng	0.00
23) Nitrobenzene-d5	7.26	82	532832	57.56	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	216913	72.53	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	1025318	59.60	ng	0.00
79) Terphenyl-d14	12.81	244	1085555	61.34	ng	0.00
Target Compounds						
49) Acenaphthylene	9.59	152	48604	2.012	ng	96
50) Dimethylphthalate	9.45	163	110733	5.861	ng	98
58) Fluorene	10.28	166	54640	3.090	ng	94
71) Phenanthrene	11.25	178	329203	13.232	ng	99
72) Anthracene	11.29	178	82349	3.240	ng	98
75) Fluoranthene	12.43	202	274350	10.220	ng	96
78) Pyrene	12.66	202	353558	14.083	ng	99
81) Benzo(a)anthracene	13.85	228	147608m	7.535	ng	
83) Chrysene	13.89	228	146233	7.346	ng	98
86) Indeno(1,2,3-cd)pyrene	16.64	276	87872	5.008	ng	97
88) Benzo(b)fluoranthene	14.87	252	153573m	8.320	ng	
89) Benzo(k)fluoranthene	14.90	252	56571m	3.552	ng	
90) Benzo(a)pyrene	15.22	252	140523	9.055	ng	100
92) Benzo(a,h,i)perylene	17.05	276	97719	7.201	ng	95

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

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