

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060520\
 Data File : BF120563.D
 Acq On : 5 Jun 2020 16:54
 Operator : JU/CG
 Sample : L2890-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26572-25062-251686

Manual Integrations
 APPROVED

mohammad
 6/8/2020 12:51:13 PM

Quant Time: Jun 06 00:35:51 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 05 15:36:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	182124	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	654695	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	302906	20.00	ng	0.00
64) Phenanthrene-d10	11.34	188	477978	20.00	ng	0.00
76) Chrysene-d12	13.98	240	375800	20.00	ng	0.00
86) Perylene-d12	15.43	264	360089	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	932374	97.56	ng	0.00
7) Phenol-d6	6.46	99	1141017	96.87	ng	0.00
23) Nitrobenzene-d5	7.38	82	751186	75.05	ng	0.00
42) 2,4,6-Tribromophenol	10.65	330	281122	100.17	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	1353527	75.82	ng	0.00
79) Terphenyl-d14	12.93	244	1138369	67.41	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.57	163	215738	11.392	ng	99
71) Phenanthrene	11.36	178	181953	8.639	ng	99
72) Anthracene	11.42	178	46761	2.213	ng	98
75) Fluoranthene	12.55	202	335527	14.806	ng	97
78) Pyrene	12.78	202	280516	10.733	ng	99
81) Benzo(a)anthracene	13.97	228	171363	8.508	ng	98
83) Chrysene	14.00	228	150068	7.107	ng	97
84) Bis(2-ethylhexyl)phthalate	13.96	149	25682	2.005	ng	# 98
87) Indeno(1,2,3-cd)pyrene	16.86	276	66007	3.351	ng	99
88) Benzo(b)fluoranthene	15.01	252	213905m	10.695	ng	
89) Benzo(k)fluoranthene	15.03	252	52629m	2.819	ng	
90) Benzo(a)pyrene	15.36	252	126602	7.237	ng	# 94
92) Benzo(a,h,i)perylene	17.29	276	61382	3.876	ng	98

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

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