

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF060120\
 Data File : BF120474.D
 Acq On : 1 Jun 2020 16:39
 Operator : JU/CG
 Sample : L2773-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM74-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/2/2020 3:26:03 PM

Quant Time: Jun 01 17:35:11 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 29 12:39:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	248692	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	903722	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	453850	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	777283	20.00	ng	0.00
76) Chrysene-d12	14.00	240	537466	20.00	ng	0.00
86) Perylene-d12	15.46	264	541637	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	1132559	86.79	ng	0.00
7) Phenol-d6	6.47	99	1485126	92.33	ng	0.00
23) Nitrobenzene-d5	7.40	82	983070	71.16	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	428433	101.88	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	1813451	67.79	ng	0.00
79) Terphenyl-d14	12.95	244	1714076	70.97	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	9.59	163	328769	11.59	ng	99
71) Phenanthrene	11.39	178	412150	12.03	ng	97
72) Anthracene	11.43	178	102387	2.98	ng	99
75) Fluoranthene	12.57	202	681971	18.51	ng	98
78) Pyrene	12.80	202	614890	16.45	ng	98
81) Benzo(a)anthracene	13.99	228	310295	10.77	ng	97
83) Chrysene	14.03	228	292806	9.70	ng	97
87) Indeno(1,2,3-cd)pyrene	16.90	276	163817m	5.53	ng	
88) Benzo(b)fluoranthene	15.03	252	391528m	13.01	ng	
89) Benzo(k)fluoranthene	15.06	252	127134m	4.53	ng	
90) Benzo(a)pyrene	15.39	252	268070	10.19	ng	96
92) Benzo(g,h,i)perylene	17.33	276	148138	6.22	ng	98

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

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