

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF060120\
 Data File : BF120472.D
 Acq On : 1 Jun 2020 15:45
 Operator : JU/CG
 Sample : L2773-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM10-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/2/2020 3:25:34 PM

Quant Time: Jun 01 16:59:17 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	225674	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	832820	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	417291	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	729400	20.00	ng	0.00
76) Chrysene-d12	14.00	240	535763	20.00	ng	0.00
86) Perylene-d12	15.46	264	554626	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	923215	77.96	ng	0.00
7) Phenol-d6	6.47	99	1208041	82.77	ng	0.00
23) Nitrobenzene-d5	7.40	82	772292	60.66	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	346229	89.55	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	1485792	60.41	ng	0.00
79) Terphenyl-d14	12.95	244	1458317	60.57	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	9.59	163	283957	10.88	ng	99
71) Phenanthrene	11.39	178	257665	8.02	ng	96
72) Anthracene	11.43	178	69735	2.16	ng	98
75) Fluoranthene	12.57	202	425236	12.30	ng	95
78) Pyrene	12.80	202	384975	10.33	ng	98
81) Benzo(a)anthracene	13.99	228	205046	7.14	ng	99
83) Chrysene	14.02	228	181042	6.01	ng	96
87) Indeno(1,2,3-cd)pyrene	16.89	276	96619	3.19	ng	95
88) Benzo(b)fluoranthene	15.03	252	251221m	8.16	ng	
89) Benzo(k)fluoranthene	15.06	252	66277m	2.30	ng	
90) Benzo(a)pyrene	15.39	252	168998	6.27	ng	98
92) Benzo(g,h,i)perylene	17.33	276	88809	3.64	ng	97

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

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