

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF060120\
 Data File : BF120475.D
 Acq On : 1 Jun 2020 17:06
 Operator : JU/CG
 Sample : L2773-16
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM61-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/2/2020 3:26:06 PM

Quant Time: Jun 02 09:24:05 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	243201	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	880831	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	418045	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	700270	20.00	ng	0.00
76) Chrysene-d12	14.00	240	520607	20.00	ng	0.00
86) Perylene-d12	15.46	264	528604	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	1075202	84.25	ng	0.00
7) Phenol-d6	6.47	99	1361480	86.56	ng	0.00
23) Nitrobenzene-d5	7.40	82	884535	65.69	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	354637	91.56	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	1599776	64.93	ng	0.00
79) Terphenyl-d14	12.95	244	1479768	63.25	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	9.59	163	259466	9.93	ng	99
71) Phenanthrene	11.39	178	161529	5.23	ng	98
75) Fluoranthene	12.57	202	422650	12.73	ng	97
78) Pyrene	12.80	202	416387	11.50	ng	96
81) Benzo(a)anthracene	13.99	228	260347m	9.33	ng	
83) Chrysene	14.03	228	251957	8.61	ng	97
87) Indeno(1,2,3-cd)pyrene	16.90	276	135097	4.67	ng	# 94
88) Benzo(b)fluoranthene	15.03	252	336083m	11.45	ng	
89) Benzo(k)fluoranthene	15.06	252	108816m	3.97	ng	
90) Benzo(a)pyrene	15.39	252	226189	8.81	ng	97
92) Benzo(g,h,i)perylene	17.34	276	121981	5.25	ng	98

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

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