

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060220\
 Data File : BF120499.D
 Acq On : 2 Jun 2020 17:12
 Operator : JU/CG
 Sample : L2773-26
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM3-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/3/2020 4:36:15 PM

Quant Time: Jun 02 19:21:31 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 02 11:37:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	146242	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	493065	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	236381	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	467293	20.00	ng	0.00
76) Chrysene-d12	14.01	240	427134	20.00	ng	0.00
86) Perylene-d12	15.47	264	391939	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	709893	92.51	ng	0.00
7) Phenol-d6	6.47	99	848904	89.75	ng	0.00
23) Nitrobenzene-d5	7.40	82	516890	68.57	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	233684	106.70	ng	-0.01
45) 2-Fluorobiphenyl	9.20	172	980485	70.38	ng	0.00
79) Terphenyl-d14	12.96	244	1212148	63.15	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	150565	10.189	ng	100
71) Phenanthrene	11.39	178	118303	5.746	ng	99
75) Fluoranthene	12.58	202	283161	12.781	ng	98
78) Pyrene	12.81	202	279156	9.398	ng	97
81) Benzo(a)anthracene	14.00	228	158634	6.930	ng	97
83) Chrysene	14.03	228	143753	5.990	ng	96
87) Indeno(1,2,3-cd)pyrene	16.93	276	77220	3.602	ng	95
88) Benzo(b)fluoranthene	15.04	252	186510m	8.568	ng	
89) Benzo(k)fluoranthene	15.07	252	62814m	3.091	ng	
90) Benzo(a)pyrene	15.41	252	126923	6.666	ng	98
92) Benzo(g,h,i)perylene	17.37	276	75740	4.394	ng	94

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

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