

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
 Data File : BF120457.D
 Acq On : 29 May 2020 23:29
 Operator : MA/SJ
 Sample : L2773-19
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM58-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/1/2020 12:21:52 PM

Quant Time: May 30 05:13:27 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	251374	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	891489	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	406626	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	638269	20.00	ng	0.00
76) Chrysene-d12	14.00	240	521404	20.00	ng	0.00
86) Perylene-d12	15.45	264	538819	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	1133278	85.92	ng	0.00
7) Phenol-d6	6.47	99	1475571	90.76	ng	0.00
23) Nitrobenzene-d5	7.40	82	963406	70.69	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	354454	94.08	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	1755701	73.26	ng	0.00
79) Terphenyl-d14	12.94	244	1556067	66.41	ng	0.00
Target Compounds						
49) Acenaphthylene	9.73	152	122231	3.596	ng	98
50) Dimethylphthalate	9.59	163	282827	11.126	ng	99
71) Phenanthrene	11.38	178	614531	21.851	ng	98
72) Anthracene	11.43	178	139116	4.930	ng	97
75) Fluoranthene	12.57	202	1284694	42.453	ng	100
78) Pyrene	12.80	202	1094591	30.187	ng	99
81) Benzo(a)anthracene	13.99	228	642234	22.983	ng	97
83) Chrysene	14.02	228	532764	18.185	ng	96
87) Indeno(1,2,3-cd)pyrene	16.89	276	291466	9.890	ng	93
88) Benzo(b)fluoranthene	15.03	252	686329m	22.933	ng	
89) Benzo(k)fluoranthene	15.06	252	238342m	8.531	ng	
90) Benzo(a)pyrene	15.39	252	486952	18.602	ng	99
91) Dibenzo(a,h)anthracene	16.89	278	72476	3.016	ng	93
92) Benzo(g,h,i)perylene	17.32	276	255406	10.778	ng	98

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

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