

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050120\
 Data File : BF120276.D
 Acq On : 1 May 2020 23:32
 Operator : MA/SJ
 Sample : L2376-03 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3

Manual Integrations
 APPROVED

mohammad
 5/4/2020 3:45:49 PM

Quant Time: May 02 03:37:34 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 01 15:26:36 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	39002	20.00	ng	0.00
21) Naphthalene-d8	7.93	136	139869	20.00	ng	0.00
39) Acenaphthene-d10	9.69	164	63517	20.00	ng	-0.01
64) Phenanthrene-d10	11.17	188	115098	20.00	ng	0.00
76) Chrysene-d12	13.81	240	100547	20.00	ng	-0.02
87) Perylene-d12	15.22	264	98048	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	129869	57.55	ng	0.00
7) Phenol-d6	6.30	99	149070	51.26	ng	-0.01
23) Nitrobenzene-d5	7.21	82	84873	31.39	ng	-0.02
42) 2,4,6-Tribromophenol	10.48	330	29726	38.22	ng	-0.01
45) 2-Fluorobiphenyl	9.01	172	192452	43.02	ng	-0.01
79) Terphenyl-d14	12.76	244	187188	31.33	ng	-0.01
Target Compounds						
49) Acenaphthylene	9.55	152	33079	5.267	ng	Qvalue 95
50) Dimethylphthalate	9.41	163	16816	3.423	ng	98
71) Phenanthrene	11.20	178	77520	12.655	ng	95
72) Anthracene	11.25	178	25118	4.014	ng	98
75) Fluoranthene	12.39	202	148690	22.496	ng	95
78) Pyrene	12.62	202	151868	17.917	ng	97
81) Benzo(a)anthracene	13.80	228	85037m	12.856	ng	
83) Chrysene	13.84	228	85758	12.760	ng	95
86) Indeno(1,2,3-cd)pyrene	16.56	276	41707	7.040	ng	96
88) Benzo(b)fluoranthene	14.82	252	97983m	13.892	ng	
89) Benzo(k)fluoranthene	14.84	252	32281m	5.304	ng	
90) Benzo(a)pyrene	15.16	252	80297	13.540	ng	98
91) Dibenzo(a,h)anthracene	16.57	278	11938	2.273	ng	# 69
92) Benzo(g,h,i)perylene	16.97	276	42873	8.268	ng	97

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

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