

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091420\
 Data File : BF121561.D
 Acq On : 14 Sep 2020 15:30
 Operator : JU/CG
 Sample : L3993-09
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PAV-B38-091120-10-17

Quant Time: Sep 14 16:52:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 10 16:47:54 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	0.00	152	0	0.00	ng	-6.83
21) Naphthalene-d8	0.00	136	0	0.00	ng	-8.12
39) Acenaphthene-d10	0.00	164	0	0.00	ng	-9.87
64) Phenanthrene-d10	0.00	188	0	0.00	ng	-11.35
76) Chrysene-d12	14.19	240	147	20.00	ng	0.20
86) Perylene-d12	15.66	264	1539	20.00	ng	0.22
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	836844	0.00	ng	0.01
7) Phenol-d6	6.47	99	1074820	0.00	ng	0.01
23) Nitrobenzene-d5	7.39	82	704681	0.00	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	308353	0.00	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	1254340	0.00	ng	0.00
79) Terphenyl-d14	12.93	244	1251651	172498.18	ng	0.00
Target Compounds						
77) Benzidine	12.93	184	17326	3555.256	ng	Qvalue # 90
78) Pyrene	12.79	202	61960	5768.945	ng	98
80) Butylbenzylphthalate	13.56	149	165	37.986	ng	# 55
81) Benzo(a)anthracene	14.01	228	29998	3225.609	ng	98
82) 3,3'-Dichlorobenzidine	14.15	252	215	59.217	ng	# 1
83) Chrysene	14.01	228	29998	3185.375	ng	96
84) Bis(2-ethylhexyl)phthalate	13.97	149	14184	2444.316	ng	96
85) Di-n-octyl phthalate	14.80	149	2899	283.081	ng	# 1
87) Indeno(1,2,3-cd)pyrene	17.30	276	17603	175.636	ng	95
88) Benzo(b)fluoranthene	15.32	252	17382	168.948	ng	# 64
89) Benzo(k)fluoranthene	15.32	252	17382	184.493	ng	# 64
90) Benzo(a)pyrene	15.47	252	6849	78.315	ng	# 9
91) Dibenzo(a,h)anthracene	17.10	278	2339	28.035	ng	# 16
92) Benzo(a,h,i)perylene	17.30	276	16118	196.538	ng	96

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

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