

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012720\
 Data File : BF118722.D
 Acq On : 27 Jan 2020 16:41
 Operator : CG/JU
 Sample : L1263-04 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CTY-SB-0203-0-11

Manual Integrations
 APPROVED

mohammad
 1/28/2020 1:13:14 PM

Quant Time: Jan 27 17:50:38 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 27 14:24:10 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	331433	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1186985	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	587420	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	899869	20.00	ng	0.00
76) Chrysene-d12	14.02	240	611957	20.00	ng	0.00
87) Perylene-d12	15.49	264	760643	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	879737	49.30	ng	0.00
7) Phenol-d6	6.48	99	1061766	45.78	ng	-0.01
23) Nitrobenzene-d5	7.42	82	653031	30.77	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	282413	41.57	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	1279272	35.79	ng	0.00
79) Terphenyl-d14	12.96	244	1082340	38.58	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.60	163	151735	3.965	ng	99
58) Fluorene	10.44	166	75293	2.156	ng	98
71) Phenanthrene	11.40	178	700096	16.126	ng	97
72) Anthracene	11.46	178	193759	4.511	ng	96
75) Fluoranthene	12.59	202	736688	16.089	ng	97
78) Pyrene	12.82	202	672595	16.189	ng	97
81) Benzo(a)anthracene	14.00	228	343941	9.343	ng	99
83) Chrysene	14.04	228	326471	9.253	ng	98
84) Bis(2-ethylhexyl)phthalate	14.00	149	42666	2.054	ng	98
86) Indeno(1,2,3-cd)pyrene	16.95	276	190853	6.226	ng	94
88) Benzo(b)fluoranthene	15.06	252	440319m	9.209	ng	
89) Benzo(k)fluoranthene	15.08	252	130506m	2.946	ng	
90) Benzo(a)pyrene	15.42	252	307231	7.449	ng	97
92) Benzo(a,h,i)perylene	17.39	276	175665	4.959	ng	97

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

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