

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042420\
 Data File : BF120202.D
 Acq On : 24 Apr 2020 15:12
 Operator : MA/SJ
 Sample : L2378-07
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3-S

Manual Integrations
 APPROVED

mohammad
 4/27/2020 3:04:41 PM

Quant Time: Apr 24 16:12:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 24 10:24:30 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	166475	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	608044	20.00	ng	0.00
39) Acenaphthene-d10	9.70	164	281306	20.00	ng	0.00
64) Phenanthrene-d10	11.19	188	428286	20.00	ng	0.00
76) Chrysene-d12	13.82	240	298054	20.00	ng	-0.01
87) Perylene-d12	15.22	264	259329	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	825732	85.72	ng	0.00
7) Phenol-d6	6.32	99	1057989	85.24	ng	0.00
23) Nitrobenzene-d5	7.23	82	657328	55.93	ng	0.00
42) 2,4,6-Tribromophenol	10.49	330	207404	60.22	ng	0.00
45) 2-Fluorobiphenyl	9.03	172	1229820	62.07	ng	0.00
79) Terphenyl-d14	12.77	244	941971	53.18	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	9.42	163	176328	8.104	ng	100
71) Phenanthrene	11.20	178	154644	6.784	ng	96
75) Fluoranthene	12.39	202	204861	8.330	ng	95
78) Pyrene	12.62	202	179591	7.147	ng	95
81) Benzo(a)anthracene	13.81	228	97920	4.994	ng	94
83) Chrysene	13.84	228	84400	4.236	ng	95
88) Benzo(b)fluoranthene	14.83	252	103668m	5.557	ng	
90) Benzo(a)pyrene	15.16	252	67918	4.330	ng	100
92) Benzo(g,h,i)perylene	16.97	276	34240	2.497	ng	# 89

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

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