

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062220\
 Data File : BF120696.D
 Acq On : 23 Jun 2020 1:02
 Operator : JU/CG
 Sample : L3079-02
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JFK-SB-IB-1.0-5.0

Manual Integrations
 APPROVED

Sohil
 6/23/2020 2:03:49 PM

Quant Time: Jun 23 04:19:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 11 11:28:22 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	147865	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	556800	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	282219	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	551179	20.00	ng	0.00
76) Chrysene-d12	13.96	240	426874	20.00	ng	0.00
86) Perylene-d12	15.42	264	451033	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	708900	82.19	ng	0.00
7) Phenol-d6	6.43	99	878529	81.78	ng	0.00
23) Nitrobenzene-d5	7.36	82	561264	56.34	ng	-0.01
42) 2,4,6-Tribromophenol	10.63	330	283687	86.46	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1105465	59.93	ng	0.00
79) Terphenyl-d14	12.91	244	1318421	68.22	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.55	163	177039	8.710	ng	98
75) Fluoranthene	12.53	202	160590	5.126	ng	98
78) Pyrene	12.76	202	158185	5.231	ng	98
81) Benzo(a)anthracene	13.95	228	91007	3.549	ng	97
83) Chrysene	13.99	228	74687	2.839	ng	96
87) Indeno(1,2,3-cd)pyrene	16.84	276	55290	2.015	ng	97
88) Benzo(b)fluoranthene	14.99	252	109753m	4.085	ng	
90) Benzo(a)pyrene	15.35	252	81499	3.389	ng	# 91
92) Benzo(g,h,i)perylene	17.27	276	51815	2.337	ng	# 93

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

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