

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
 Data File : BF120454.D
 Acq On : 29 May 2020 22:07
 Operator : MA/SJ
 Sample : L2773-23
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM89-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/1/2020 12:21:41 PM

Quant Time: May 30 04:54:11 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 29 12:39:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	286177	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	1026301	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	473732	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	737150	20.00	ng	0.00
76) Chrysene-d12	13.99	240	563788	20.00	ng	0.00
86) Perylene-d12	15.44	264	572853	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	1549569	103.19	ng	0.00
7) Phenol-d6	6.47	99	1941687	104.91	ng	0.00
23) Nitrobenzene-d5	7.40	82	1295539	82.57	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	468354	106.70	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	2185440	78.27	ng	0.00
79) Terphenyl-d14	12.95	244	1884127	74.37	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	419533	14.166	ng	99
71) Phenanthrene	11.38	178	86646	2.668	ng	97
75) Fluoranthene	12.57	202	194965	5.578	ng	95
78) Pyrene	12.79	202	179717	4.584	ng	98
81) Benzo(a)anthracene	13.98	228	101300	3.353	ng	95
83) Chrysene	14.02	228	99223	3.132	ng	95
88) Benzo(b)fluoranthene	15.02	252	134815m	4.237	ng	
90) Benzo(a)pyrene	15.38	252	92567	3.326	ng	98
92) Benzo(g,h,i)perylene	17.31	276	56750	2.253	ng	97

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

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