

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF052820\
 Data File : BF120428.D
 Acq On : 29 May 2020 00:46
 Operator : MA/SJ
 Sample : L2744-20 2X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM50-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 5/29/2020 10:29:17 PM

Quant Time: May 29 05:54:48 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 28 13:47:54 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	255861	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	875389	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	383734	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	641946	20.00	ng	0.00
76) Chrysene-d12	14.00	240	554526	20.00	ng	0.00
86) Perylene-d12	15.45	264	468883	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	555491	41.38	ng	0.00
7) Phenol-d6	6.46	99	671188	40.56	ng	0.00
23) Nitrobenzene-d5	7.40	82	393762	29.42	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	141561	39.82	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	776960	34.35	ng	0.00
79) Terphenyl-d14	12.94	244	734418	29.47	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	117541	4.90	ng	99
71) Phenanthrene	11.38	178	104588	3.70	ng	98
75) Fluoranthene	12.57	202	238172	7.83	ng	98
78) Pyrene	12.80	202	241077	6.25	ng	99
81) Benzo(a)anthracene	13.99	228	146075	4.92	ng	98
83) Chrysene	14.02	228	142044m	4.56	ng	
87) Indeno(1,2,3-cd)pyrene	16.88	276	58344	2.28	ng	96
88) Benzo(b)fluoranthene	15.03	252	159694m	6.13	ng	
90) Benzo(a)pyrene	15.39	252	105790	4.64	ng	98
92) Benzo(g,h,i)perylene	17.32	276	55009	2.67	ng	97

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

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