

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060320\
 Data File : BF120517.D
 Acq On : 3 Jun 2020 18:47
 Operator : JU/CG
 Sample : L2839-09
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM102-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/4/2020 12:12:40 PM

Quant Time: Jun 04 00:56:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 03 14:26:17 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	131628	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	480511	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	237032	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	492386	20.00	ng	0.00
76) Chrysene-d12	14.00	240	354988	20.00	ng	-0.01
86) Perylene-d12	15.45	264	296057	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	659026	95.42	ng	0.00
7) Phenol-d6	6.47	99	832128	97.75	ng	0.00
23) Nitrobenzene-d5	7.40	82	555079	75.56	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	254392	115.83	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	1074003	76.88	ng	0.00
79) Terphenyl-d14	12.95	244	1160273	72.73	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	204173	13.778	ng	99
71) Phenanthrene	11.38	178	161025	7.422	ng	96
75) Fluoranthene	12.57	202	260500	11.159	ng	99
78) Pyrene	12.80	202	249371	10.101	ng	97
81) Benzo(a)anthracene	13.99	228	133993	7.043	ng	100
83) Chrysene	14.02	228	128785m	6.457	ng	
87) Indeno(1,2,3-cd)pyrene	16.90	276	59691	3.686	ng	95
88) Benzo(b)fluoranthene	15.03	252	147373m	8.962	ng	
89) Benzo(k)fluoranthene	15.05	252	48000m	3.127	ng	
90) Benzo(a)pyrene	15.39	252	87979	6.117	ng	97
92) Benzo(a,h,i)perylene	17.33	276	57355	4.405	ng	100

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

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