

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM051019\
 Data File : BM020262.D
 Acq On : 11 May 2019 05:06
 Operator : JU/SJ
 Sample : K2779-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RT-5124

Manual Integrations
 APPROVED

mohammad
 5/13/2019 8:25:36 AM

Quant Time: May 11 06:18:31 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM043019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 03:58:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	100123	20.00	ng	-0.02
21) Naphthalene-d8	10.57	136	415118	20.00	ng	-0.02
39) Acenaphthene-d10	14.42	164	283721	20.00	ng	-0.02
64) Phenanthrene-d10	17.17	188	660059	20.00	ng	-0.01
76) Chrysene-d12	21.35	240	505111	20.00	ng	-0.02
87) Perylene-d12	23.64	264	497338	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	695541	113.55	ng	-0.01
7) Phenol-d6	6.95	99	1048481	125.30	ng	0.00
23) Nitrobenzene-d5	8.94	82	712785	67.79	ng	-0.02
42) 2,4,6-Tribromophenol	15.92	330	539859	122.59	ng	0.00
45) 2-Fluorobiphenyl	13.04	172	1579469	68.49	ng	-0.02
79) Terphenyl-d14	19.80	244	2353111	81.26	ng	-0.01
Target Compounds				Qvalue		
50) Dimethylphthalate	13.87	163	285772	11.851	ng	98
75) Fluoranthene	19.23	202	172362	3.739	ng	98
78) Pyrene	19.59	202	153693	4.532	ng	100
81) Benzo(a)anthracene	21.34	228	112051	3.163	ng	98
83) Chrysene	21.39	228	112545	3.276	ng	97
86) Indeno(1,2,3-cd)pyrene	25.96	276	84581	2.070	ng	# 100
88) Benzo(b)fluoranthene	22.94	252	178566m	5.437	ng	
90) Benzo(a)pyrene	23.54	252	113902	3.818	ng	# 86
92) Benzo(g,h,i)perylene	26.68	276	86012	2.920	ng	# 94

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

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