

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
 Data File : BM020297.D
 Acq On : 12 May 2019 14:20
 Operator : JU/SJ
 Sample : K2766-09
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P026-SS007-1824-01

Manual Integrations
 APPROVED

Sohil
 5/14/2019 8:07:28 PM

Quant Time: May 13 06:53:25 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	7.76	152	55888	20.00	ng	-0.03
21) Naphthalene-d8	10.56	136	190882	20.00	ng	-0.03
39) Acenaphthene-d10	14.41	164	116929	20.00	ng	-0.02
64) Phenanthrene-d10	17.16	188	299819	20.00	ng	-0.02
76) Chrysene-d12	21.34	240	340393	20.00	ng	-0.02
87) Perylene-d12	23.63	264	405955	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	405503	118.60	ng	-0.02
7) Phenol-d6	6.93	99	532207	113.94	ng	-0.02
23) Nitrobenzene-d5	8.93	82	357001	73.84	ng	-0.03
42) 2,4,6-Tribromophenol	15.90	330	230675	127.10	ng	-0.02
45) 2-Fluorobiphenyl	13.03	172	710737	74.79	ng	-0.03
79) Terphenyl-d14	19.79	244	1372480	70.33	ng	-0.02
Target Compounds						
49) Acenaphthylene	14.13	152	40979	3.332	ng	Qvalue 97
50) Dimethylphthalate	13.87	163	58706	5.907	ng	97
71) Phenanthrene	17.20	178	141925	8.575	ng	98
75) Fluoranthene	19.22	202	271935	12.986	ng	99
78) Pyrene	19.59	202	531211	23.244	ng	100
81) Benzo(a)anthracene	21.33	228	203773m	8.536	ng	
83) Chrysene	21.39	228	256257m	11.069	ng	
86) Indeno(1,2,3-cd)pyrene	25.95	276	140467	5.101	ng	# 100
88) Benzo(b)fluoranthene	22.94	252	264971m	9.884	ng	
89) Benzo(k)fluoranthene	22.97	252	74103m	3.030	ng	
90) Benzo(a)pyrene	23.53	252	182804	7.507	ng	97
92) Benzo(a,h,i)perylene	26.66	276	169125	7.035	ng	98

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

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