

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050719\
 Data File : BM020186.D
 Acq On : 08 May 2019 04:00
 Operator : JU/SJ
 Sample : K2672-20
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-10-COMP

Manual Integrations
 APPROVED

Sohil
 5/9/2019 6:44:38 PM

Quant Time: May 08 07:03:20 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	7.79	152	140423	20.00	ng	0.00
21) Naphthalene-d8	10.59	136	485255	20.00	ng	0.00
39) Acenaphthene-d10	14.44	164	290740	20.00	ng	0.00
64) Phenanthrene-d10	17.19	188	680129	20.00	ng	0.00
76) Chrysene-d12	21.37	240	731646	20.00	ng	0.00
87) Perylene-d12	23.66	264	873623	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	1482638	172.59	ng	0.00
7) Phenol-d6	6.97	99	2040920	173.90	ng	0.02
23) Nitrobenzene-d5	8.96	82	1369738	111.44	ng	0.00
42) 2,4,6-Tribromophenol	15.93	330	905929	200.75	ng	0.01
45) 2-Fluorobiphenyl	13.06	172	2320623	98.21	ng	0.00
79) Terphenyl-d14	19.81	244	3771016	89.90	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.89	163	394248	15.955	ng	99
71) Phenanthrene	17.23	178	232804	6.201	ng	99
75) Fluoranthene	19.24	202	527553	11.106	ng	98
78) Pyrene	19.61	202	502066	10.221	ng	98
81) Benzo(a)anthracene	21.35	228	324852	6.331	ng	92
83) Chrysene	21.40	228	301273m	6.054	ng	
86) Indeno(1,2,3-cd)pyrene	26.00	276	285576	4.825	ng	# 100
88) Benzo(b)fluoranthene	22.97	252	530657m	9.199	ng	
89) Benzo(k)fluoranthene	23.01	252	148262m	2.817	ng	
90) Benzo(a)pyrene	23.56	252	381665	7.284	ng	# 87
92) Benzo(g,h,i)perylene	26.72	276	286069m	5.529	ng	

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

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