

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050619\
 Data File : BM020156.D
 Acq On : 07 May 2019 06:26
 Operator : JU/SJ
 Sample : K2673-06 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-13-COMP

Manual Integrations
 APPROVED

mohammad
 5/7/2019 3:12:12 PM

Quant Time: May 07 07:42:23 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	153460	20.00	ng	0.00
21) Naphthalene-d8	10.59	136	554287	20.00	ng	0.00
39) Acenaphthene-d10	14.44	164	311620	20.00	ng	0.00
64) Phenanthrene-d10	17.19	188	718459	20.00	ng	0.00
76) Chrysene-d12	21.37	240	741114	20.00	ng	0.00
87) Perylene-d12	23.67	264	862357	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	291476	31.05	ng	0.00
7) Phenol-d6	6.96	99	400933	31.26	ng	0.00
23) Nitrobenzene-d5	8.96	82	261690	18.64	ng	0.00
42) 2,4,6-Tribromophenol	15.93	330	143943	29.76	ng	0.01
45) 2-Fluorobiphenyl	13.05	172	445799	17.60	ng	0.00
79) Terphenyl-d14	19.81	244	642394	15.12	ng	0.00
Target Compounds						
49) Acenaphthylene	14.16	152	76900	2.346	ng	Qvalue 97
50) Dimethylphthalate	13.89	163	72254	2.728	ng	98
71) Phenanthrene	17.23	178	526903	13.286	ng	99
72) Anthracene	17.32	178	203908	5.142	ng	98
75) Fluoranthene	19.24	202	1370062	27.303	ng	99
78) Pyrene	19.61	202	1135548	22.822	ng	99
81) Benzo(a)anthracene	21.36	228	753508	14.498	ng	96
83) Chrysene	21.41	228	593156m	11.768	ng	
86) Indeno(1,2,3-cd)pyrene	26.00	276	565580	9.433	ng	# 100
88) Benzo(b)fluoranthene	22.97	252	952768	16.731	ng	# 96
89) Benzo(k)fluoranthene	23.01	252	330865m	6.370	ng	
90) Benzo(a)pyrene	23.57	252	797033	15.409	ng	# 95
91) Dibenzo(a,h)anthracene	26.01	278	152668	2.838	ng	# 81
92) Benzo(g,h,i)perylene	26.73	276	512898	10.043	ng	96

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

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