

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050619\
 Data File : BM020157.D
 Acq On : 07 May 2019 07:02
 Operator : JU/SJ
 Sample : K2673-15 5X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP-01-COMP

Manual Integrations
 APPROVED

mohammad
 5/7/2019 3:12:15 PM

Quant Time: May 07 07:44:57 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	155726	20.00	ng	0.00
21) Naphthalene-d8	10.59	136	550478	20.00	ng	0.00
39) Acenaphthene-d10	14.44	164	311062	20.00	ng	0.00
64) Phenanthrene-d10	17.19	188	727404	20.00	ng	0.00
76) Chrysene-d12	21.37	240	755665	20.00	ng	0.00
87) Perylene-d12	23.67	264	876498	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	165062	17.33	ng	0.00
7) Phenol-d6	6.96	99	231524	17.79	ng	0.00
23) Nitrobenzene-d5	8.96	82	143330	10.28	ng	0.00
42) 2,4,6-Tribromophenol	15.93	330	84661	17.53	ng	0.00
45) 2-Fluorobiphenyl	13.05	172	299355	11.84	ng	0.00
79) Terphenyl-d14	19.81	244	466189	10.76	ng	0.00
Target Compounds						
71) Phenanthrene	17.23	178	576428	14.356	ng	99
72) Anthracene	17.32	178	176220	4.389	ng	99
75) Fluoranthene	19.24	202	1370403	26.974	ng	99
78) Pyrene	19.61	202	1392541	27.448	ng	99
81) Benzo(a)anthracene	21.36	228	803625	15.164	ng	98
83) Chrysene	21.41	228	654776	12.740	ng	96
86) Indeno(1,2,3-cd)pyrene	26.00	276	492686	8.059	ng	# 100
88) Benzo(b)fluoranthene	22.97	252	942420m	16.283	ng	
89) Benzo(k)fluoranthene	23.01	252	315731m	5.980	ng	
90) Benzo(a)pyrene	23.57	252	737518	14.028	ng	# 94
91) Dibenzo(a,h)anthracene	26.00	278	130958	2.395	ng	# 79
92) Benzo(a,h,i)perylene	26.73	276	495059	9.537	ng	98

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

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