

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050619\
 Data File : BM020136.D
 Acq On : 06 May 2019 15:24
 Operator : JU/SJ
 Sample : K2642-01 2X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OK-02-050219

Manual Integrations
 APPROVED

mohammad
 5/7/2019 3:11:39 PM

Quant Time: May 07 05:04:06 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	135887	20.00	ng	0.00
21) Naphthalene-d8	10.59	136	507674	20.00	ng	0.00
39) Acenaphthene-d10	14.43	164	306712	20.00	ng	0.00
64) Phenanthrene-d10	17.18	188	728473	20.00	ng	0.00
76) Chrysene-d12	21.37	240	741485	20.00	ng	0.00
87) Perylene-d12	23.66	264	884310	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	383445	46.12	ng	0.00
7) Phenol-d6	6.95	99	549271	48.36	ng	0.00
23) Nitrobenzene-d5	8.95	82	372056	28.93	ng	0.00
42) 2,4,6-Tribromophenol	15.93	330	234388	49.23	ng	0.00
45) 2-Fluorobiphenyl	13.05	172	795151	31.90	ng	0.00
79) Terphenyl-d14	19.81	244	1289480	30.33	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.89	163	105144	4.033	ng	Qvalue # 96
71) Phenanthrene	17.22	178	293847	7.307	ng	99
72) Anthracene	17.32	178	98260	2.444	ng	99
75) Fluoranthene	19.24	202	551196	10.834	ng	99
78) Pyrene	19.60	202	514527	10.335	ng	99
81) Benzo(a)anthracene	21.35	228	284292	5.467	ng	95
83) Chrysene	21.40	228	264006m	5.235	ng	
86) Indeno(1,2,3-cd)pyrene	26.00	276	171019	2.851	ng	# 100
88) Benzo(b)fluoranthene	22.97	252	329805m	5.648	ng	
89) Benzo(k)fluoranthene	23.01	252	127571m	2.395	ng	
90) Benzo(a)pyrene	23.56	252	256056	4.827	ng	# 82
92) Benzo(a,h,i)perylene	26.71	276	174561	3.333	ng	# 82

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

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