

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060319\
 Data File : BG041047.D
 Acq On : 3 Jun 2019 23:13
 Operator : HP/JU
 Sample : K2641-19
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CL-02-053119-D

Manual Integrations
 APPROVED

mohammad
 6/5/2019 8:35:30 AM

Quant Time: Jun 04 09:54:18 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 29 18:39:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	86280	20.00	ng	-0.01
21) Naphthalene-d8	10.94	136	320882	20.00	ng	0.00
39) Acenaphthene-d10	14.75	164	213729	20.00	ng	-0.01
64) Phenanthrene-d10	17.50	188	471954	20.00	ng	0.00
76) Chrysene-d12	21.77	240	465698	20.00	ng	-0.01
87) Perylene-d12	25.12	264	511985	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	470743	98.38	ng	0.00
7) Phenol-d6	7.27	99	681214	92.19	ng	0.00
23) Nitrobenzene-d5	9.30	82	482620	66.77	ng	0.00
42) 2,4,6-Tribromophenol	16.24	330	294480	92.81	ng	0.00
45) 2-Fluorobiphenyl	13.37	172	902211	60.79	ng	-0.01
79) Terphenyl-d14	20.09	244	1256108	57.25	ng	-0.01
Target Compounds						
50) Dimethylphthalate	14.19	163	87893	4.858	ng	97
71) Phenanthrene	17.54	178	290179	11.804	ng	99
72) Anthracene	17.62	178	51579	2.127	ng	96
75) Fluoranthene	19.54	202	463686	15.271	ng	99
78) Pyrene	19.90	202	545473	18.018	ng	98
81) Benzo(a)anthracene	21.75	228	264993	8.548	ng	97
83) Chrysene	21.82	228	288154	9.713	ng	97
86) Indeno(1,2,3-cd)pyrene	28.95	276	132284	3.640	ng	# 98
88) Benzo(b)fluoranthene	24.05	252	313004	10.079	ng	# 93
89) Benzo(k)fluoranthene	24.11	252	110999m	3.612	ng	
90) Benzo(a)pyrene	24.96	252	254668	8.596	ng	97
92) Benzo(a,h,i)perylene	30.16	276	141627	4.924	ng	# 94

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

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