

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050819\
 Data File : BM020202.D
 Acq On : 08 May 2019 19:33
 Operator : JU/SJ
 Sample : K2641-05 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CL-02-050719-C

Manual Integrations
 APPROVED

Sohil
 5/9/2019 6:45:33 PM

Quant Time: May 09 02:29:42 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	161828	20.00	ng	0.00
21) Naphthalene-d8	10.58	136	583874	20.00	ng	0.00
39) Acenaphthene-d10	14.43	164	345531	20.00	ng	0.00
64) Phenanthrene-d10	17.17	188	749051	20.00	ng	0.00
76) Chrysene-d12	21.36	240	760722	20.00	ng	0.00
87) Perylene-d12	23.66	264	886941	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	543825	54.93	ng	0.00
7) Phenol-d6	6.95	99	744480	55.04	ng	0.00
23) Nitrobenzene-d5	8.95	82	527482	35.67	ng	0.00
42) 2,4,6-Tribromophenol	15.92	330	284494	53.04	ng	0.00
45) 2-Fluorobiphenyl	13.04	172	983578	35.02	ng	-0.01
79) Terphenyl-d14	19.80	244	1381157	31.67	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.89	163	136572	4.650	ng	99
58) Fluorene	15.48	166	72474	2.510	ng	99
71) Phenanthrene	17.22	178	581502	14.064	ng	98
72) Anthracene	17.31	178	215246	5.207	ng	98
75) Fluoranthene	19.23	202	949900	18.157	ng	98
78) Pyrene	19.60	202	882466	17.278	ng	98
81) Benzo(a)anthracene	21.34	228	514264	9.640	ng	98
83) Chrysene	21.40	228	435265	8.413	ng	96
86) Indeno(1,2,3-cd)pyrene	25.99	276	288631	4.690	ng	# 100
88) Benzo(b)fluoranthene	22.96	252	593813	10.139	ng	# 96
89) Benzo(k)fluoranthene	23.00	252	200644m	3.756	ng	
90) Benzo(a)pyrene	23.55	252	455749	8.567	ng	95
92) Benzo(a,h,i)perylene	26.71	276	279407	5.319	ng	93

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

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